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May 2, 2016 **VIA PRIVATE CARRIER**

Ms. Ellen Jackson
Maryland Department of the Environment
Oil Control Program
1800 Washington Blvd.
Baltimore, MD 21230

Subject: Transmittal of the Block F Underground Storage Tank Closure Report
(Appendix H to the Construction Completion Report)
Lockheed Martin Corporation; Middle River Complex
2323 Eastern Boulevard, Middle River, Baltimore County, Maryland

Dear Ms. Jackson:

Please find enclosed two hard copies with a CD of the above-referenced document. This document presents detailing underground storage tank (UST) removal activities at the Middle River Complex in Middle River, Maryland.

We respectfully request to receive Maryland Department of the Environment's approval by May 30, 2016.

Please let me know if you have any questions. My office phone is (301) 548-2209.

Sincerely,

Thomas D. Blackman
Project Lead, Environmental Remediation

cc: (via email without enclosure)
James Carroll, MDE
Gary Schold, MDE
Mark Mank, MDE
Christine Kline, Lockheed Martin
Norman Varney, Lockheed Martin
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Doug Mettee, Lockheed Martin MST
Jann Richardson, Lockheed Martin
Justin Tetlow, MRAS

# Block F Underground Storage Tank Closure Report
# Lockheed Martin Middle River Complex
# 2323 Eastern Boulevard
# Middle River, Maryland

Facility ID – 13198

Case # 16-0084BA

Prepared for:

Lockheed Martin Corporation

Prepared by:

Tetra Tech, Inc.

May 2016

Michael Martin, P.G.
Regional Manager

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# ACRONYMS

| | |
|---|---|
| Block | Tax Block |
| ESA | environmental site assessment |
| Lockheed Martin | Lockheed Martin Corporation |
| MDE | Maryland Department of the Environment |
| µg/kg | microgram(s) per kilogram |
| MRC | Middle River Complex |
| OCP | Oil Control Program |
| PAH(s) | polycyclic aromatic hydrocarbon(s) |
| REC(s) | recognized environmental condition(s) |
| RAP | remedial action plan |
| Tetra Tech | Tetra Tech, Inc. |
| TPH | total petroleum hydrocarbons |
| UCL | upper confidence level |
| USEPA | United States Environmental Protection Agency |
| UST | underground storage tank |

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# Section 1
# Narrative

On behalf of Lockheed Martin Corporation (Lockheed Martin), Tetra Tech Inc. (Tetra Tech) has prepared this report detailing underground storage tank (UST) removal activities at the Middle River Complex (MRC), located in Baltimore County at 2323 Eastern Boulevard, Middle River, Maryland. The MRC site is comprised of approximately 160 acres with 12 main buildings. The property includes an active industrial area and yard, perimeter parking lots, an athletic field, a vacant concrete-covered lot, a trailer and parts storage lot, and numerous grassy areas along the facility perimeter. Locked chain-link fences surround all exterior lots and the main industrial area. The site is bounded by Eastern Boulevard (Route 150) to the north, Dark Head Cove to the south, Cow Pen Creek to the west, and Martin State Airport (MSA) to the east.

A Phase I environmental site assessment (ESA) was conducted at Middle River Complex in February 2003 to identify recognized environmental conditions (RECs) through a desktop study of historical documents and cursory site inspection. The Phase I ESA identified Former Aviation Fuel underground storage tanks, including seven closed-in-place 1,000-gallon underground storage tanks beneath the Boat Launch Area's concrete tarmac, as REC #5. The tanks were reportedly last used in 1954 and closed in place in 1986-1987. During the Phase I ESA investigation, no closure records were available for these underground storage tanks, and MRC personnel interviewed did not know whether confirmatory sampling had been conducted as part of the tank closure.

Due to the unknown status of the underground storage tanks, Lockheed Martin included tank removal and disposal into the Block F soil remedial action plan (RAP) which evaluated soil remedial alternatives for the remediation of polycyclic aromatic hydrocarbons (PAHs), expressed as benzo(a)pyrene equivalent concentrations, in Block F soil (Tetra Tech, 2013). The Block F soil remedial action plan was submitted and approved by Maryland Department of the Environment Controlled Hazardous Substance Enforcement Division in December 2013. The underground storage tank removal was then incorporated in the *100% Soil Remedial Action Design for Blocks D,*

*D Panhandle, F, G, and H* (Tetra Tech, 2015b) that was approved by state regulators in January 2015.

The seven 1,000-gallon underground storage tanks and a network of piping were uncovered approximately seven feet below ground surface in the concrete tarmac area of Block F during remedial activities on August 3-10, 2015 (Appendix A). All seven tanks had been previously filled with a concrete slurry. A sheet piling system was installed to stabilize the excavation area and safely remove the tanks.

The seven 1,000-gallon tanks were removed on November 4, 2015 and confirmation samples were collected from the base of the excavation on November 5, 2015 as directed by Ellen Jackson, the onsite Maryland Department of the Environment Oil Control Program (OCP) inspector. No sidewall samples were collected from the underground storage tank area because the sidewalls were covered by the sheet-piling system installed to stabilize the excavation area during tank removal. No evidence of contamination (via visual inspection or elevated photoionization detector readings) were encountered during excavation and removal of the tanks. The tanks were cut, cleaned, and secured for recycling at an approved recycling facility.

The piping network extended approximately 100-feet from the area where the underground storage tanks were removed. The piping network consisted of seven two-inch steel pipes and two three-inch steel pipes, located approximately six feet below ground surface. All piping within the erosion and sediment controls was removed on November 13, 2015. The remaining pipe located outside of the erosion and sediment control was removed on December 28, 2015 when no rain was forecasted and the excavation area was backfilled on the same day. Confirmation base and sidewall samples were collected from the pipe excavation areas as directed by the Oil Control Program inspector. No visually impacted material or elevated photoionization detector readings above background concentrations were encountered during excavation and removal of the piping network. The pipes were cut, cleaned, and secured for off-site disposal at an approved waste disposal facility.

Photographic documentation of the removal of the underground storage tank and associated piping is in Appendix B. The subsequent sections of this closure report detail the soil sampling methodology, sampling results, and recommendations.

# Section 2
# Soil Sampling

Fourteen soil samples (F-VB-UST-1 through F-VB-UST-14) were collected from the bottom of the underground storage tank (UST) excavation area following removal of the tanks at Tax Block (Block) F. Verification samples were collected from a depth of fourteen feet below ground surface in areas directed by the onsite Maryland Department of the Environment Oil Control Program (OCP) inspector. No sidewall samples were collected from the UST area because of the sheet piling system installed to stabilize the excavation area.

Additionally, eight soil samples (F-VB-UST-P-1 through F-VB-UST-P-8) were collected from the bottom of the pipe excavation area and four soil samples (F-VS-UST-P-1 through F-VS-UST-P-4) were collected from the sidewalls of the pipe excavation area. Pipe verification base samples were collected as grab samples from the excavation base at a frequency of one sample per fifteen linear feet. Pipe verification sidewall samples were collected as grab samples at locations directed by the OCP inspector, at the depth the pipes were before removal.

All soil samples were placed in wide-mouth laboratory-supplied glass jars and immediately placed on ice in a clean cooler. The cooler temperature was maintained at less than or equal to four degrees Celsius. Soil samples were delivered to TestAmerica Inc. and analyzed as directed by the onsite OCP inspector for full-suite volatile organic compounds by United States Environmental Protection Agency (USEPA) Method 8260, for semivolatile organic compounds by USEPA Method 8270, for total petroleum hydrocarbon (TPH) -diesel-range organics and TPH-gasoline-range organics by Method 8015, and for polychlorinated biphenyls by USEPA method 8082.

Per the OCP inspector’s guidance, collection of groundwater samples was not required after tanks were removed. Sampling locations are in Appendix C.

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# Section 3
# Analytical Results

Table 3-1 presents positive detections in the fourteen soil samples (F-VB-UST-1 through F-VB-UST-14) collected after the underground storage tanks (USTs) were removed. The soil samples were compared to Maryland Department of the Environment (MDE) soil standards (MDE, 2008). No exceedances of non-residential screening criteria were detected in the soil samples collected beneath the seven USTs. The complete analytical results for the base samples are in Appendix E.

Table 3-2 presents the detections in the soil sidewall and base samples collected after the piping network was removed. The pipe soil samples were compared to MDE soil standards. One exceedance of the MDE non-residential cleanup standard for benzo(a)pyrene (390 micrograms per kilogram [µg/kg]) was detected in the pipe removal area in sample F-VS-UST-P-3 at a concentration of 700 µg/kg. As presented in Section 1, polycyclic aromatic hydrocarbons, expressed as benzo(a)pyrene equivalents (BaPEq)[1], are the contaminant of concern evaluated in the Tax Block F soil remedial action plan (RAP). The BaPEq 95% upper confidence limit (UCL) for all pipe verification base and sidewall samples was calculated and compared to the site-specific remedial goal of 2,900 µg/kg as defined in the Block F RAP developed for the protection of industrial workers. As discussed below, the BaPEq RG for the industrial worker is also protective of the construction worker or utility worker (the likely receptor for subsurface soils).

The BaPEq 95% UCL was calculated to be 810 µg/kg which is less than the industrial worker remedial goal of 2,900 µg/kg. Additionally, the cancer risk calculated for the future industrial worker was $3\times10^{-6}$, a value less than the MDE cancer risk limit of $1x10^{-5}$. The BaPEq 95% UCL

[1] Per current United States Environmental Protection Agency (USEPA) guidance (USEPA, 1993b), concentrations of benzo(a)pyrene, benzo(a)anthracene, benzo(b)fluoranthene, benzo(k)fluoranthene, chrysene, dibenz(a,h)anthracene, and indeno(1,2,3-c,d)pyrene can be combined in an adjusted risk-weighted concentration that is expressed in terms of benzo(a)pyrene toxicity. This calculated value, called a *benzo(a)pyrene equivalent concentration* (BaPEq), uses USEPA-recommended toxicity equivalency factors (TEFs) to estimate the potency of each of these polycyclic aromatic hydrocarbon (PAH) compounds relative to that of benzo(a)pyrene. The TEFs are then used to convert each individual PAH concentration into an equivalent concentration of benzo(a)pyrene; these values are summed to arrive at the calculated BaPEq concentration.

of 810 µg/kg was also less than a construction worker remedial goal of 16,000 µg/kg. It should be noted that remedial goals developed for BaPEq for construction workers or utility workers exceed those developed for a typical industrial worker primarily because the exposure frequency/duration assumptions for the typical industrial worker are more conservative than those for the construction worker or utility worker. Thus, the BaPEq remedial goal for the industrial worker is also protective of the construction worker or utility worker. The remedial goal of 16,000 µg/kg, based on a target risk level of $1\times10^{-5}$, was developed for construction workers exposed to BaPEq in subsurface soil in the 2012 human health risk assessment (HHRA) (Tetra Tech, 2012) (Appendix D).

A residual risk technical memorandum detailing the UCL calculation along with the ProUCL calculation output sheet and risk is in Appendix D[2]. Laboratory analytical reports are in Appendix E.

[2] Note that the subsurface soil samples collected for the Block F UST closure are not included in the surface soil data set used for the residual risk analysis (RRA) conducted for the remaining Block F soils (see Appendix D). The RRA evaluates cancer and noncancer risks to industrial workers exposed to surface (the top two feet) soil.

**Table 3-1**

**Summary of Soil Verification Samples - Block F Underground Storage Tank**
**Block F Underground Storage Tank Closure Report**
**Lockheed Martin Corporation Middle River Complex**
**Middle River, Maryland**


| LOCATION<br>SAMPLE ID<br>SAMPLE DATE | Remedial Goal | MDE Residential Cleanup Standard | MDE Non-Residential Cleanup Standard | F-VB-UST-1<br>F-VB-UST-1-14<br>20151105 | F-VB-UST-2<br>F-VB-UST-2-14<br>20151105 | F-VB-UST-2<br>F-VB-UST-2-14-AVG<br>20151105 | F-VB-UST-2<br>F-VB-UST-2-14-D<br>20151105 | F-VB-UST-3<br>F-VB-UST-3-14<br>20151105 | F-VB-UST-4<br>F-VB-UST-4-14<br>20151105 |
|---|---|---|---|---|---|---|---|---|---|
| **VOLATILES (UG/KG)** | | | | | | | | | |
| ACETONE | NA | 7,000,000 | 92,000,000 | 5.7 U | 6.3 U | 6.15 U | 6 UJ | 6.6 U | 6.2 U |
| NAPHTHALENE | NA | 160,000 | 2,000,000 | 0.35 U | 0.38 U | 0.375 U | 0.37 U | 0.41 U | 0.38 U |
| TOLUENE | NA | 630,000 | 8,200,000 | 0.27 U | 0.3 U | 0.29 U | 0.28 U | 0.31 U | 0.29 U |
| **SEMIVOLATILES (UG/KG)** | | | | | | | | | |
| ACETOPHENONE | NA | NA | NA | 16 J | 11 U | 11 U | 11 U | 11 U | 13 J |
| BENZALDEHYDE | NA | NA | NA | 14 UJ | 14 UJ | 14 U | 14 UJ | 15 UJ | 14 UJ |
| BIS(2-ETHYLHEXYL)PHTHALATE | NA | 46,000 | 200,000 | 31 J | 22 U | 16.5 | 22 J | 23 U | 33 J |
| DI-N-BUTYL PHTHALATE | NA | 780,000 | 10,000,000 | 18 U | 18 U | 17.5 U | 17 U | 18 U | 18 U |
| BAP EQUIVALENT-HALFND | 2900 | NA | NA | 0.76 U | 0.75 U | 0.75 U | 0.75 U | 0.78 U | 0.76 U |
| BAP EQUIVALENT-POS | 2900 | NA | NA | 0.76 U | 0.75 U | 0.75 U | 0.75 U | 0.78 U | 0.76 U |
| **PETROLEUM HYDROCARBONS** | | | | | | | | | |
| TPH (C10-C32) (mg/kg) | NA | 230 | 620 | 16 U | 10 U | 9.1 U | 8.2 U | 7.7 U | 7.6 U |
| TPH (C06-C10) (ug/kg) | NA | 230,000 | 620,000 | 55 U | 54 U | 54 U | 54 U | 56 U | 55 U |

NOTES:

No sample exceeded the RG or the MDE residential and non-residential cleanup standards

J - Estimated
MDE - Maryland Department of the Environmen
NA - Not applicable/not available
U - Non-detect

**Table 3-1**

**Summary of Soil Verification Samples - Block F Underground Storage Tank**
**Block F Underground Storage Tank Closure Report**
**Lockheed Martin Corporation Middle River Complex**
**Middle River, Maryland**


| LOCATION<br>SAMPLE ID<br>SAMPLE DATE | **Remedial Goal** | **MDE Residential Cleanup Standard** | **MDE Non-Residential Cleanup Standard** | **F-VB-UST-5<br>F-VB-UST-5-14<br>20151105** | **F-VB-UST-6<br>F-VB-UST-6-14<br>20151105** | **F-VB-UST-7<br>F-VB-UST-7-14<br>20151105** | **F-VB-UST-8<br>F-VB-UST-8-14<br>20151105** | **F-VB-UST-9<br>F-VB-UST-9-14<br>20151105** | **F-VB-UST-10<br>F-VB-UST-10-14<br>20151105** |
|---|---|---|---|---|---|---|---|---|---|
| **VOLATILES (UG/KG)** | | | | | | | | | |
| ACETONE | NA | 7,000,000 | 92,000,000 | 13 J | 6.6 U | 6.4 U | 6.6 U | 12 J | 6.4 U |
| NAPHTHALENE | NA | 160,000 | 2,000,000 | 0.41 J | 0.45 J | 0.39 U | 0.4 U | 0.4 U | 0.39 U |
| TOLUENE | NA | 630,000 | 8,200,000 | 0.31 U | 0.31 U | 0.3 U | 0.33 J | 0.36 J | 0.3 U |
| **SEMIVOLATILES (UG/KG)** | | | | | | | | | |
| ACETOPHENONE | NA | NA | NA | 11 U | 11 U | 11 U | 11 U | 13 J | 11 U |
| BENZALDEHYDE | NA | NA | NA | 14 UJ | 14 UJ | 15 UJ | 15 UJ | 15 UJ | 14 UJ |
| BIS(2-ETHYLHEXYL)PHTHALATE | NA | 46,000 | 200,000 | 38 J | 32 J | 32 J | 23 U | 40 J | 29 J |
| DI-N-BUTYL PHTHALATE | NA | 780,000 | 10,000,000 | 18 U | 18 U | 18 U | 18 U | 19 U | 17 U |
| BAP EQUIVALENT-HALFND | 2900 | NA | NA | 0.75 U | 0.76 U | 0.78 U | 0.78 U | 0.8 U | 0.74 U |
| BAP EQUIVALENT-POS | 2900 | NA | NA | 0.75 U | 0.76 U | 0.78 U | 0.78 U | 0.8 U | 0.74 U |
| **PETROLEUM HYDROCARBONS** | | | | | | | | | |
| TPH (C10-C32) (mg/kg) | NA | 230 | 620 | 7.5 U | 7.7 U | 16 U | 7.7 U | 7.8 U | 13 U |
| TPH (C06-C10) (ug/kg) | NA | 230,000 | 620,000 | 77 J | 55 U | 55 U | 56 U | 57 U | 53 U |

NOTES:
No sample exceeded the RG or the MDE residential and non-residential cleanup standards

J - Estimated
MDE - Maryland Department of the Environmen
NA - Not applicable/not available
U - Non-detect

**Table 3-1**

**Summary of Soil Verification Samples - Block F Underground Storage Tank**
**Block F Underground Storage Tank Closure Report**
**Lockheed Martin Corporation Middle River Complex**
**Middle River, Maryland**


| LOCATION<br>SAMPLE ID<br>SAMPLE DATE | **Remedial Goal** | **MDE Residential Cleanup Standard** | **MDE Non-Residential Cleanup Standard** | **F-VB-UST-11<br>F-VB-UST-11-14<br>20151105** | **F-VB-UST-12<br>F-VB-UST-12-14<br>20151105** | **F-VB-UST-13<br>F-VB-UST-13-14<br>20151105** | **F-VB-UST-14<br>F-VB-UST-14-14<br>20151105** |
|---|---|---|---|---|---|---|---|
| **VOLATILES (UG/KG)** | | | | | | | |
| ACETONE | NA | 7,000,000 | 92,000,000 | 6.8 U | 6.3 U | 5.9 U | 6 U |
| NAPHTHALENE | NA | 160,000 | 2,000,000 | 0.42 U | 0.39 U | 0.36 U | 0.37 U |
| TOLUENE | NA | 630,000 | 8,200,000 | 0.32 U | 0.3 J | 0.28 U | 0.29 J |
| **SEMIVOLATILES (UG/KG)** | | | | | | | |
| ACETOPHENONE | NA | NA | NA | 38 J | 11 U | 11 U | 11 J |
| BENZALDEHYDE | NA | NA | NA | 19 J | 14 UJ | 14 UJ | 14 UJ |
| BIS(2-ETHYLHEXYL)PHTHALATE | NA | 46,000 | 200,000 | 23 U | 23 U | 60 J | 22 U |
| DI-N-BUTYL PHTHALATE | NA | 780,000 | 10,000,000 | 18 U | 18 U | 18 J | 18 U |
| BAP EQUIVALENT-HALFND | 2900 | NA | NA | 0.77 U | 0.77 U | 0.74 U | 0.76 U |
| BAP EQUIVALENT-POS | 2900 | NA | NA | 0.77 U | 0.77 U | 0.74 U | 0.76 U |
| **PETROLEUM HYDROCARBONS** | | | | | | | |
| TPH (C10-C32) (mg/kg) | NA | 230 | 620 | 9 U | 30 U | 13 U | 11 U |
| TPH (C06-C10) (ug/kg) | NA | 230,000 | 620,000 | 55 U | 54 U | 54 U | 54 U |

NOTES:

No sample exceeded the RG or the MDE residential and non-residential cleanup standards

J - Estimated
MDE - Maryland Department of the Environmen
NA - Not applicable/not available
U - Non-detect

**Table 3-2**

**Summary of Soil Verification Samples - Block F Underground Storage Tank Pipe**
**Block F Underground Storage Tank Closure Report**
**Lockheed Martin Corporation Middle River Complex**
**Middle River, Maryland**


| LOCATION | Remedial Goal | MDE Residential Cleanup Standard | MDE Non-Residential Cleanup Standard | F-VB-UST-P-1 | F-VB-UST-P-2 | F-VB-UST-P-3 | F-VB-UST-P-4 | F-VB-UST-P-5 | F-VB-UST-P-6 | F-VB-UST-P-7 | F-VB-UST-P-8 | F-VS-UST-P-1 | F-VS-UST-P-1 | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| SAMPLE ID | | | | | | | | | | | | | F-VS-UST-P-1-AVG | F-VS-UST-P-1-D | F-VS-UST-P-2 | F-VS-UST-P-3 | F-VS-UST-P-4 |
| SAMPLE DATE | | | | 20151113 | 20151113 | 20151113 | 20151113 | 20151113 | 20151113 | 20151228 | 20151228 | 20151113 | 20151113 | 20151113 | 20151113 | 20151113 | 20151113 |
| **VOLATILES (UG/KG)** | | | | | | | | | | | | | | | | | |
| 1,2,3-TRIMETHYLBENZENE | NA | NA | NA | 0.54 U | 0.51 U | 0.54 U | 0.49 U | 0.5 U | 0.53 U | 0.86 J | 0.5 U | 0.87 J | 0.555 | 0.48 U | 98 | 1.5 J | 0.53 UJ |
| 1,2,4-TRIMETHYLBENZENE | NA | NA | NA | 0.68 U | 0.65 U | 0.68 U | 0.62 U | 0.63 U | 0.67 U | 1.9 J | 0.64 U | 0.66 U | 0.63 U | 0.6 U | 170 | 1.6 J | 0.67 UJ |
| 2-BUTANONE | NA | 4,700,000 | 61,000,000 | 1.3 J | 5.4 J | 4.9 J | 1.9 J | 2 J | 4.1 J | 4.8 U | 1.9 U | 18 J | 12.2 | 6.4 J | 8.8 J | 2.4 J | 4.4 J |
| 4-ISOPROPYLTOLUENE | NA | NA | NA | 0.48 U | 0.46 U | 0.48 U | 0.44 U | 0.44 U | 0.47 U | 0.48 U | 0.57 J | 0.47 U | 0.445 U | 0.42 U | 6.5 | 0.44 U | 0.47 UJ |
| ACETONE | NA | 7,000,000 | 92,000,000 | 8.2 J | 70 | 66 | 16 J | 11 J | 57 | 36 | 9.4 U | 86 J | 60 | 34 J | 26 U | 12 J | 21 J |
| BENZENE | NA | 12,000 | 52,000 | 1.1 U | 1 U | 1.1 U | 0.97 U | 0.97 U | 1 U | 1 U | 0.99 U | 1.8 J | 1.1325 | 0.93 U | 6.8 | 0.96 U | 1 U |
| ETHYLBENZENE | NA | 780,000 | 10,000,000 | 0.37 U | 0.36 U | 0.37 U | 0.34 U | 0.35 U | 0.37 U | 0.66 J | 0.35 U | 0.36 U | 0.345 U | 0.33 U | 140 | 0.34 U | 0.37 U |
| ISOPROPYLBENZENE | NA | 780,000 | 10,000,000 | 0.23 U | 0.22 U | 0.23 U | 0.21 U | 0.22 U | 0.23 U | 0.24 J | 2 J | 14 | 8.4 | 2.8 J | 63 | 0.21 U | 0.23 U |
| M+P-XYLENES | NA | NA | NA | 0.49 U | 0.47 U | 0.49 U | 0.45 U | 1.4 J | 0.49 U | 2.9 J | 0.46 U | 1.9 J | 1.0575 | 0.43 U | 520 J | 0.45 U | 0.48 U |
| NAPHTHALENE | NA | 160,000 | 2,000,000 | 0.41 U | 0.39 U | 0.41 U | 0.38 U | 0.38 U | 0.4 U | 0.61 U | 0.88 U | 1.1 J | 0.64 | 0.36 U | 17 | 1.7 J | 0.88 J |
| N-BUTYLBENZENE | NA | NA | NA | 0.69 U | 0.66 U | 0.69 U | 0.63 U | 0.64 U | 0.68 U | 0.69 U | 0.83 J | 7.1 | 4.4 | 1.7 J | 11 | 0.63 U | 0.68 UJ |
| N-PROPYLBENZENE | NA | NA | NA | 0.7 ` | 0.67 U | 0.7 U | 0.64 U | 0.65 U | 0.69 U | 0.7 U | 1.9 J | 20 J | 12.2 | 4.4 J | 55 | 0.64 U | 0.69 UJ |
| O-XYLENE | NA | NA | NA | 0.67 U | 0.64 U | 0.67 U | 0.61 U | 0.62 U | 0.66 U | 1.2 J | 0.63 U | 0.75 J | 0.5225 | 0.59 U | 14 | 0.61 U | 0.65 U |
| SEC-BUTYLBENZENE | NA | NA | NA | 1.1 U | 1 U | 1.1 U | 0.98 U | 0.98 U | 1.1 U | 1.1 U | 1 U | 8.6 | 5.3 | 2 J | 12 | 0.97 U | 1 UJ |
| TERT-BUTYLBENZENE | NA | NA | NA | 0.78 U | 0.75 U | 0.78 U | 0.72 U | 0.72 U | 0.77 U | 0.78 U | 0.74 U | 2.3 J | 1.3225 | 0.69 U | 4.4 J | 0.71 U | 0.77 UJ |
| TOLUENE | NA | 630,000 | 8,200,000 | ND | ND | ND | ND | ND | ND | 1.6 J | 0.3 U | ND | ND | ND | ND | ND | ND |
| TOTAL XYLENES | NA | 1,600,000 | 20,000,000 | 0.63 U | 0.6 U | 0.63 U | 0.58 U | 1.4 J | 0.62 U | 4.1 J | 0.59 U | 2.7 J | 1.49 | 0.56 U | 530 | 0.57 U | 0.62 U |
| **SEMIVOLATILES (UG/KG)** | | | | | | | | | | | | | | | | | |
| 1,1-BIPHENYL | NA | NA | NA | 4.1 U | 4.1 U | 4.3 U | 4.2 U | 4.2 U | 4.1 U | 4.3 U | 4.1 U | 4.2 U | 4.15 U | 4.1 U | 7.2 J | 85 J | 48 J |
| 2-METHYLNAPHTHALENE | NA | 31,000 | 410,000 | 0.59 U | 0.59 U | 0.61 U | 0.6 U | 0.61 U | 0.59 U | 0.61 U | 53 | 7.1 J | 7.45 | 7.8 | 85 | 360 | 220 |
| ACENAPHTHENE | NA | 470,000 | 6,100,000 | 0.89 U | 0.9 U | 0.93 U | 0.91 U | 0.92 U | 0.89 U | 0.93 U | 63 | 0.91 U | 0.9 U | 0.89 U | 10 | 350 | 110 |
| ACENAPHTHYLENE | NA | 470,000 | 6,100,000 | 0.41 U | 0.41 U | 0.43 U | 0.42 U | 0.42 U | 0.41 U | 0.43 U | 0.41 U | 0.42 U | 2.105 | 4 J | 4.6 J | 28 | 23 |
| ACETOPHENONE | NA | NA | NA | 11 U | 11 U | 11 U | 11 U | 11 U | 11 U | 11 U | 11 U | 11 U | 11.25 | 17 J | 24 J | 27 U | 11 U |
| ANTHRACENE | NA | 2,300,000 | 31,000,000 | 0.91 U | 0.92 U | 0.96 U | 0.94 U | 0.95 U | 0.92 U | 0.95 U | 97 | 4.6 J | 5 | 5.4 J | 7.4 J | 590 | 160 |
| BENZALDEHYDE | NA | NA | NA | 14 U | 14 U | 45 J | 14 U | 15 U | 14 U | 14 U | 15 U | 17 J | 21 | 25 J | 23 J | 35 U | 17 J |
| BENZO(A)ANTHRACENE | NA | 220 | 3,900 | 0.74 U | 0.74 U | 0.77 U | 0.76 U | 0.76 U | 0.74 U | 0.77 U | 140 | 12 | 14 | 16 | 14 | 960 | 200 |
| BENZO(A)PYRENE | NA | 22 | 390 | 0.75 U | 0.76 U | 0.78 U | 0.77 U | 0.78 U | 0.75 U | 0.78 U | 100 | 12 | 14.5 | 17 | 12 | 700 | 100 |
| BAP EQUIVALENT-HALFND | 2900 | NA | NA | 0.75 U | 0.76 U | 0.78 U | 0.77 U | 0.78 U | 0.75 U | 0.78 U | 153 | 16 | 20 | 23 | 17 | 1,005 | 168 |
| BAP EQUIVALENT-POS | 2900 | NA | NA | 0.75 U | 0.76 U | 0.78 U | 0.77 U | 0.78 U | 0.75 U | 0.78 U | 153 | 16 | 19 | 23 | 16 | 1,005 | 168 |
| BENZO(B)FLUORANTHENE | NA | 220 | 3,900 | 0.69 U | 0.7 U | 0.72 U | 0.71 U | 0.72 U | 0.69 U | 0.72 U | 180 | 17 J | 23.5 | 30 J | 17 | 1,100 | 240 |
| BENZO(G,H,I)PERYLENE | NA | 230,000 | 3,100,000 | 0.41 U | 0.41 U | 0.43 U | 0.42 U | 0.42 U | 0.41 U | 0.43 U | 51 | 8.7 | 9.35 | 10 | 10 | 240 | 41 |
| BENZO(K)FLUORANTHENE | NA | 2,200 | 39,000 | 0.8 U | 0.8 U | 0.83 U | 0.82 U | 0.82 U | 0.8 U | 0.83 U | 47 | 9.7 | 9.8 | 9.9 | 8.9 | 440 | 71 |
| BIS(2-ETHYLHEXYL)PHTHALATE | NA | 46,000 | 200,000 | 25 U | 38 U | 55 U | 30 U | 62 U | 31 U | 160 U | 180 U | 23 U | 22.5 U | 22 U | 31 U | 56 U | 33 U |
| CARBAZOLE | NA | 32,000 | 140,000 | 32 U | 32 U | 33 U | 32 U | 33 U | 32 U | 33 U | 32 U | 32 U | 32 U | 32 U | 32 U | 150 | 32 U |
| CHRYSENE | NA | 22,000 | 390,000 | 1.3 U | 1.3 U | 1.3 U | 1.3 U | 1.3 U | 1.3 U | 1.3 U | 94 | 18 | 20 | 22 | 20 | 1,000 | 260 |
| DIBENZO(A,H)ANTHRACENE | NA | 22 | 390 | 0.77 U | 0.78 U | 0.81 U | 0.79 U | 0.8 U | 0.77 U | 0.81 U | 16 | 0.79 U | 0.78 U | 0.77 U | 0.79 U | 69 | 19 |
| DIBENZOFURAN | NA | 7,800 | 100,000 | 0.77 U | 0.78 U | 0.81 U | 0.79 U | 0.8 U | 0.77 U | 0.81 U | 51 J | 14 J | 15 | 16 J | 11 J | 330 | 170 |
| FLUORANTHENE | NA | 310,000 | 4,100,000 | 0.64 U | 0.65 U | 0.67 U | 0.66 U | 0.67 U | 0.65 U | 9.3 | 490 | 27 | 32 | 37 | 30 | 2,500 | 560 |
| FLUORENE | NA | 310,000 | 4,100,000 | 0.62 U | 0.63 U | 0.65 U | 0.64 U | 0.64 U | 0.62 U | 0.65 U | 88 | 4.8 J | 5.45 | 6.1 J | 14 | 350 | 120 |
| INDENO(1,2,3-CD)PYRENE | NA | 220 | 3,900 | 0.41 U | 0.41 U | 0.43 U | 0.42 U | 0.42 U | 0.41 U | 0.43 U | 45 | 7.3 J | 8.25 | 9.2 | 9 | 250 | 40 |
| NAPHTHALENE | NA | 160,000 | 2,000,000 | 0.96 U | 0.97 U | 1 U | 0.98 U | 0.99 U | 0.96 U | 4.3 J | 14 | 7.4 J | 9.7 | 12 | 51 | 460 | 290 |
| PHENANTHRENE | NA | 2,300,000 | 31,000,000 | 0.86 U | 0.86 U | 0.9 U | 0.88 U | 0.88 U | 0.86 U | 9.3 | 570 | 20 | 23.5 | 27 | 35 | 1,900 | 550 |
| PYRENE | NA | 230,000 | 3,100,000 | 0.52 U | 0.52 U | 0.54 U | 0.53 U | 0.53 U | 0.52 U | 7.9 J | 370 | 23 | 26.5 | 30 | 28 | 1,800 | 420 |
| **PCBS (UG/KG)** | | | | | | | | | | | | | | | | | |
| AROCLOR-1260 | NA | 320 | 1400 | 10 U | 11 U | 11 U | 11 U | 11 U | 11 U | 11 U | 11 U | 11 U | 11 U | 11 U | 10 U | 11 U | 11 U |
| TOTAL AROCLOR | NA | NA | NA | 0 U | 0 U | 0 U | 0 U | 0 U | 0 U | 0 U | 0 U | 0 U | 0 U | 0 U | 0 U | 0 U | 0 U |
| TOTAL AROCLOR HALFND | NA | NA | NA | 52.15 | 53.75 | 55.95 | 53.8 | 54.35 | 53.75 | 55.95 | 53.8 | 53.75 | 53.75 | 53.75 | 52.15 | 53.8 | 53.75 |
| **PETROLEUM HYDROCARBONS (UG/KG)** | | | | | | | | | | | | | | | | | |
| TPH (C06-C10) | NA | 230000 | 620000 | 54 U | 55 U | 60 J | 55 U | 55 U | 55 U | 56 U | 55 U | 3500 J | 2650 | 1800 J | 5500 | 110 J | 55 U |
| **PETROLEUM HYDROCARBONS (MG/KG)** | | | | | | | | | | | | | | | | | |
| TPH (C10-C32) | NA | 230 | 620 | 7.6 U | 7.6 U | 7.9 U | 7.6 U | 7.8 U | 8 J | 7.9 U | 42 | 80 | 65.5 | 51 | 130 | 42 | 120 |

NOTES:
**GREY SHADED CELLS EXCEED MDE RESIDENTIAL CLEANUP STANDARDS**
BLACK SHADED CELLS EXCEED MDE RESIDENTIAL AND NON-RESIDENTIAL CLEANUP STANDARDS
No sample exceeded the remedial goal

U - Non-detect
NA - Not applicable/not available
J - Estimated

# Section 4
# Conclusions and Recommendations

Tax Block (Block F) of Middle River Complex (MRC) is an industrial site at 2323 Eastern Boulevard in Middle River, Maryland. Block F occupies approximately 11.94 acres in the southern portion of MRC. Block F is bounded on the north by the concrete foundation of former Building D (Block E) and properties formerly owned by Lockheed Martin now occupied by the Tilley Chemical Company and North American Electric. Block F is bounded to the east by the waterfront lot portion of Block D and Dark Head Cove, to the south by Dark Head Cove, and to the west by Cow Pen Creek. East of MRC is the Martin State Airport, totaling approximately 750 acres, which in 1975 was transferred from Martin Marietta to the State of Maryland.

Sampling results from the soil sample F-VS-UST-P-3 indicate the presence of benzo(a)pyrene above the Maryland Department of the Environment non-residential soil cleanup standard (390 µg/kg); however, when included in the overall benzo(a)pyrene equivalent (BaPEq) calculation of 810 µg/kg, the concentration is below the remedial goal of 2,900 µg/kg for industrial workers established and approved by the Maryland Department of the Environment for Block F (Appendix D). Additionally, the cancer risk to the industrial worker was calculated to be $3\times10^{-6}$ which is less than the $1\times10^{-5}$ cancer risk set in the remedial action plan for BaPEq in Block F soils. The BaPEq exposure point concentration calculation was also less than a remedial goal of 16,000 µg/kg based on a target risk level of $1\times10^{-5}$ developed for construction workers exposed to BaPEq in subsurface soil.

Because the current and future use of the site is industrial, and sampling results indicate the cancer risk posed by remaining benzo(a)pyrene in subsurface soil is below $1\times10^{-5}$ for industrial workers and construction workers, Lockheed Martin Corporation recommends performing no additional work with regard to the underground storage tanks removed from Block F.

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# Section 5
# References


1. Earth Tech, 2003. *Phase I Environmental Site Assessment, Chesapeake Industrial Park, Middle River, Maryland, May*.

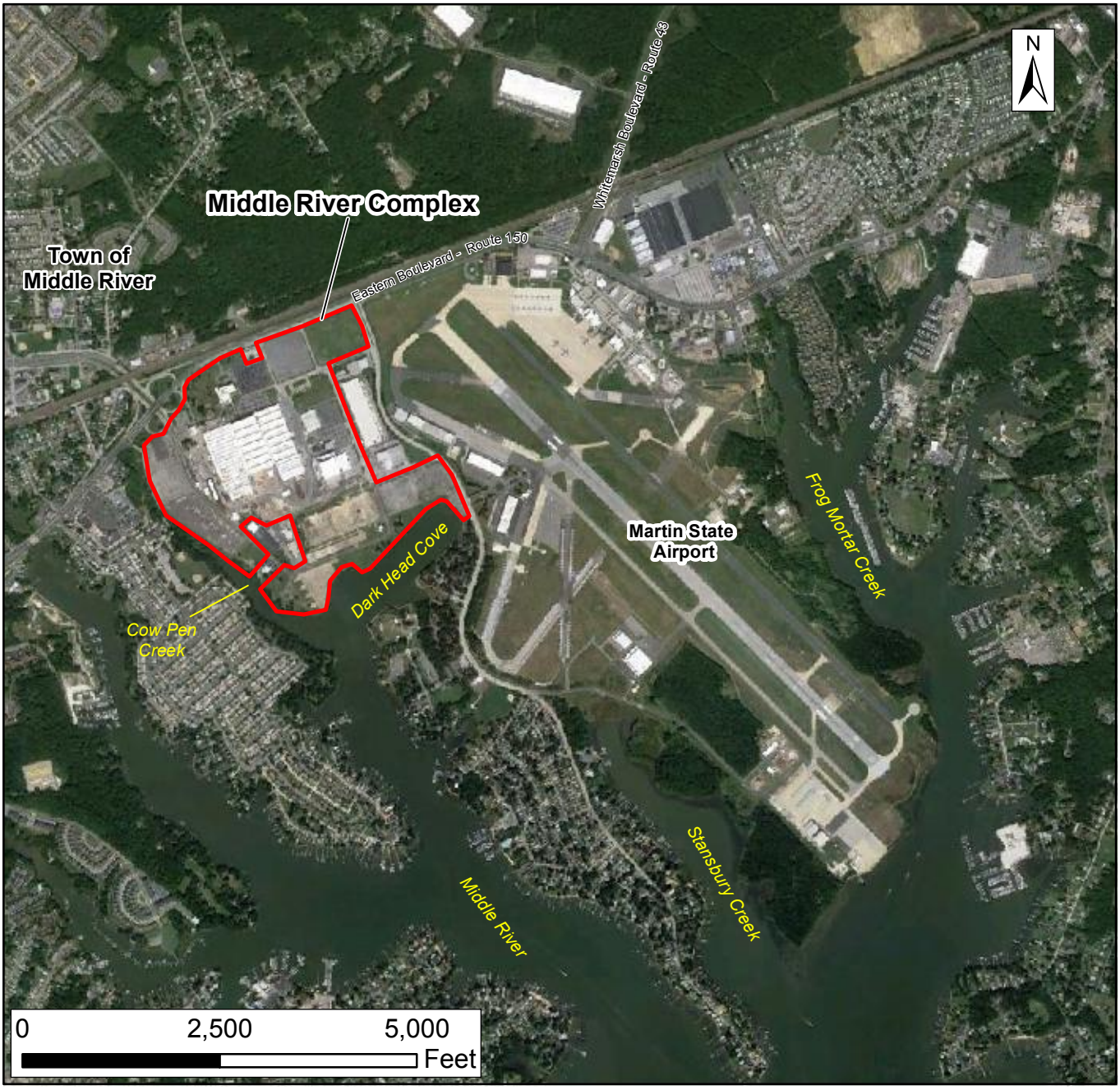
2. MDE (Maryland Department of the Environment), 2008. *Cleanup Standards for Soil and Groundwater.* Interim Final. Update No. 2.1. June.

3. Tetra Tech, Inc. (Tetra Tech), 2012. *Human Health Risk Assessment (HHRA) for Blocks D, E, F, G, and H Soils, Lockheed Martin Middle River Complex, 2323 Eastern Boulevard, Middle River, Maryland*. September.

4. Tetra Tech, Inc. (Tetra Tech), 2013. *Block F Remedial Action Plan. Lockheed Martin Middle River Complex, Middle River, Maryland*. Report prepared for Lockheed Martin Corporation, Bethesda, Maryland by Tetra Tech, Inc., Germantown, Maryland. January.

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# APPENDIX A—UST LOCATION MAP




Source: Google Earth, 2013

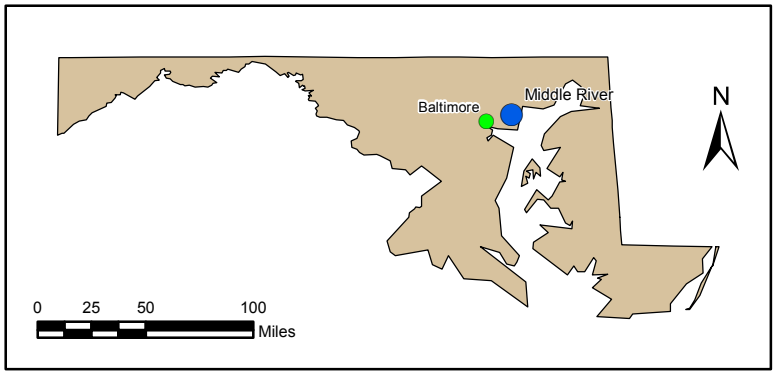



**FIGURE 1**

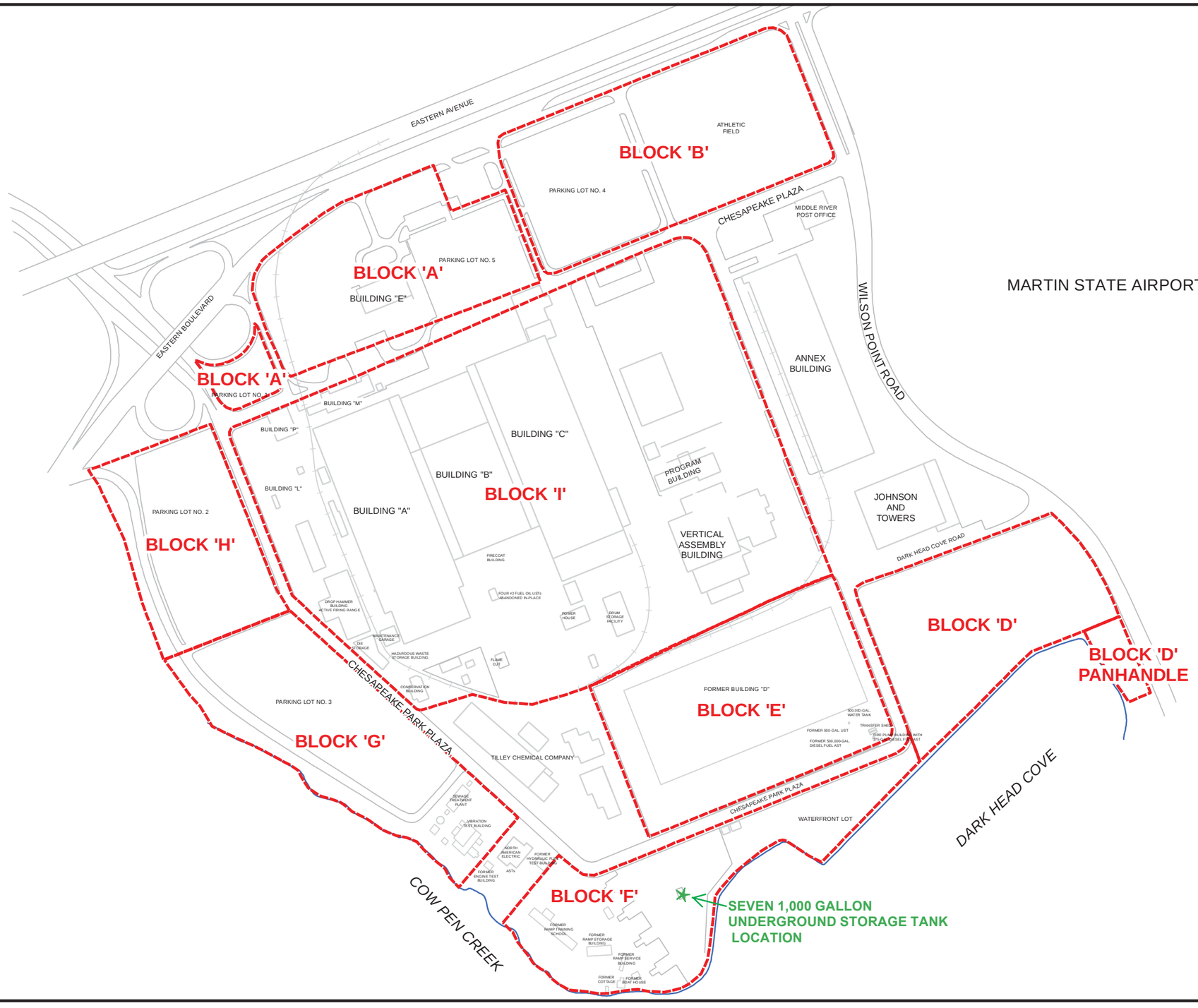
**MIDDLE RIVER COMPLEX LOCATION MAP**

***Lockheed Martin Middle River Complex***
***Middle River, Maryland***

| DATE MODIFIED: | CREATED BY: |
|---|---|
| 5/1/13 | MP |

TETRA TECH

Map Document: (K:\GProject\middle_river\Maps\MiddleRiver_MRC_Location Map_Fig 1-1_revised050113.mxd) 5/1/2013 -- 11:18:55 AM




MARTIN STATE AIRPORT

**FIGURE 2**

**LOCKHEED MARTIN**
**MIDDLE RIVER COMPLEX**
**TAX BLOCKS**

**LEGEND**

- TAX BLOCK
- STRUCTURE
- RAILROAD TRACKS

***Lockheed Martin Middle River Complex***
***Middle River, Maryland***

0 200 400 Feet

N

DATE MODIFIED: 6/20/13

CREATED BY: MP

TETRA TECH

Map Document: (K:\GProject\middle_river\Maps\MRC Tax Blocks_091911.mxd)

# APPENDIX B—PHOTOGRAPHIC DOCUMENTATION

**Photographs of Underground Storage Tank and Piping Removal**



Comments: 1. Technicians assessing tank integrity and conducting tank surveying



Comments: 2. Technician screening suction tube opening of UST #3 in Block F UST excavation.

**Photos of Select Daily Activities (continued)**



Comments: 3. Equipment operator breaks concrete retaining wall to facilitate UST removal. Note concrete between tanks.



Comments: 4. UST #7 detached from concrete retaining walls. 6"-8" concrete pad beneath USTs.

**Photos of Select Daily Activities**




Comments: 1. Block F UST area; removal of UST from Block F excavation area.




Comments: 2. Block F UST area; repositioning UST on to plastic lined area.

**Photos of Select Daily Activities (continued)**



Comments: 3. Block F UST area; USTs, concrete walls and concrete pad.



Comments: 4. Block F UST Area following removal of USTs, concrete walls and concrete pad.

**Photos of Select Daily Activities**




Comments: 1. Block F UST area; bobcat with hammer attachment utilized to break up concrete for future sampling and disposal.

Comments: 2.

**Photos of Select Daily Activities (continued)**




Comments: 3. Block F UST area; excavator utilized to dump and spread select fill during backfilling activities.




Comments: 4. Block F UST area; mini-roller utilized to compact select fill.

**Photos of Select Daily Activities**



Comments: 1. Block F UST pipe removal; (7) 2" steel pipes uncovered, the (2) 3" steel pipes run underneath the 2" pipes. Piping system located approximately 6' bgs.



Comments: 2. Block F UST pipe removal; excavator bucket removes 2" steel pipes from the excavation area.

**Photos of Select Daily Activities (continued)**




Comments: 3. Block F UST pipe removal; excavator bucket with clay soil from the base of the excavation @ 8-10' bgs, soil does not appear to be impacted and PID readings did not exceed 0.0 ppm.



Comments: 4. Block F UST pipe removal; excavator bucket with clay soil from base of excavation @ 8-10 bgs.

**Photos of Select Daily Activities**

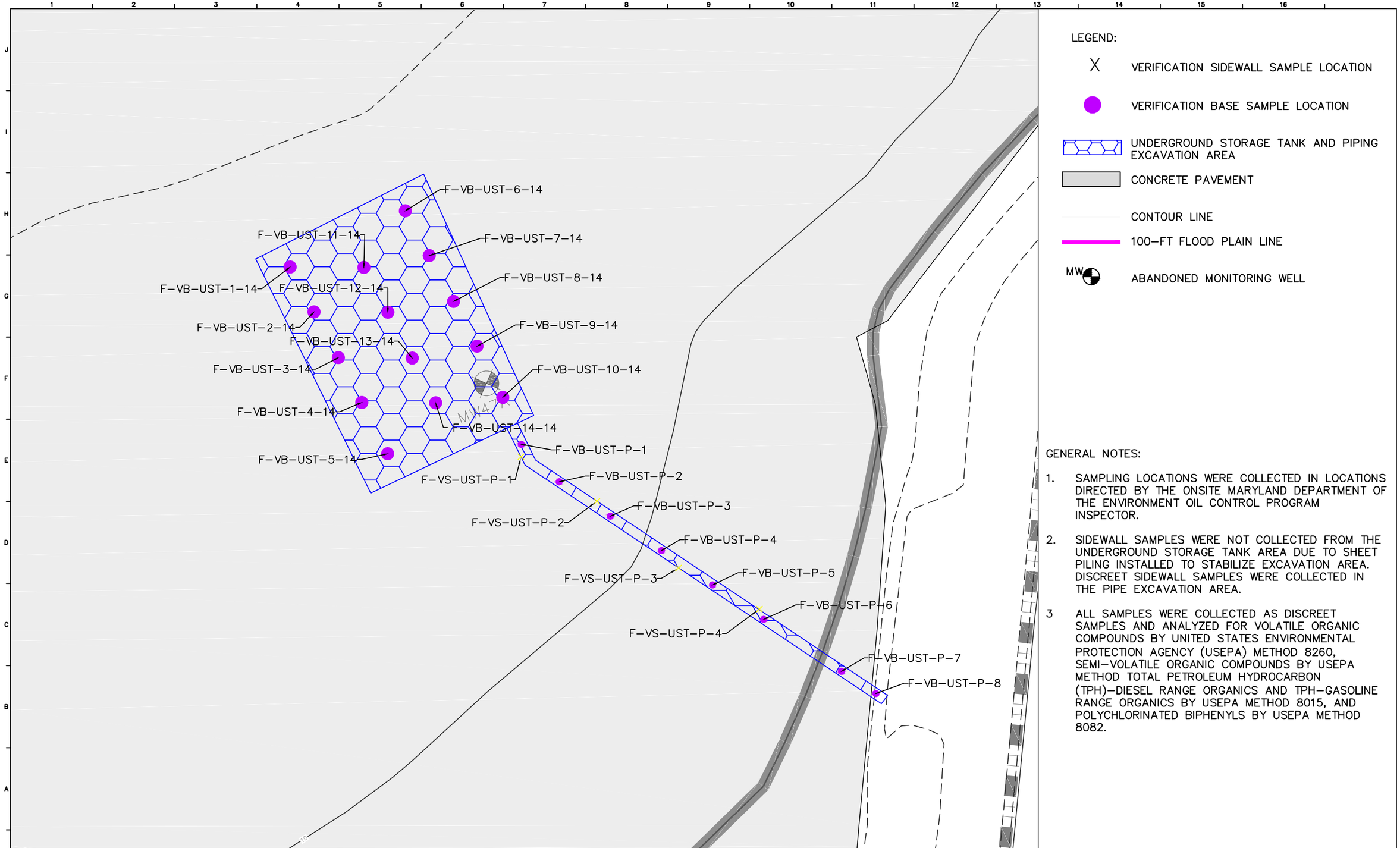


Comments: 1. Delivery of select fill to Block F UST Area prior to backfilling with CR-6 to stabilize base of excavation.



Comments: 2. Placing select fill in Block F UST excavation.

# APPENDIX C—SAMPLING LOCATION DIAGRAM




LEGEND:

- X — VERIFICATION SIDEWALL SAMPLE LOCATION
- ● — VERIFICATION BASE SAMPLE LOCATION
- UNDERGROUND STORAGE TANK AND PIPING EXCAVATION AREA
- CONCRETE PAVEMENT
- CONTOUR LINE
- 100-FT FLOOD PLAIN LINE
- MW — ABANDONED MONITORING WELL

GENERAL NOTES:

1. SAMPLING LOCATIONS WERE COLLECTED IN LOCATIONS DIRECTED BY THE ONSITE MARYLAND DEPARTMENT OF THE ENVIRONMENT OIL CONTROL PROGRAM INSPECTOR.

2. SIDEWALL SAMPLES WERE NOT COLLECTED FROM THE UNDERGROUND STORAGE TANK AREA DUE TO SHEET PILING INSTALLED TO STABILIZE EXCAVATION AREA. DISCREET SIDEWALL SAMPLES WERE COLLECTED IN THE PIPE EXCAVATION AREA.

3 ALL SAMPLES WERE COLLECTED AS DISCREET SAMPLES AND ANALYZED FOR VOLATILE ORGANIC COMPOUNDS BY UNITED STATES ENVIRONMENTAL PROTECTION AGENCY (USEPA) METHOD 8260, SEMI-VOLATILE ORGANIC COMPOUNDS BY USEPA METHOD TOTAL PETROLEUM HYDROCARBON (TPH)-DIESEL RANGE ORGANICS AND TPH-GASOLINE RANGE ORGANICS BY USEPA METHOD 8015, AND POLYCHLORINATED BIPHENYLS BY USEPA METHOD 8082.

TETRA TECH
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GERMANTOWN, MD 20874
T: (301) 528-5552 | F: (301) 528-3000

| MARK | DATE | DESCRIPTION | BY |
|---|---|---|---|
| | | | |

LOCKHEED MARTIN
MIDDLE RIVER COMPLEX

UNDERGROUND STORAGE TANK AND PIPING AREA
SAMPLING LOCATION DIAGRAM
BLOCK F

| | |
|---|---|
| DATE: | 1/22/2016 |
| PROJECT NO.: | 112IC07483 |
| DESIGNED BY: | ISG |
| DRAWN BY: | ISG |
| CHECKED BY: | SAN |
| SHEET: | 1 OF 1 |

COPYRIGHT TETRA TECH INC.

FIGURE 3

K:\CADD\ARC\112IC07483\Construction Completion\Block F\AS BUILT FIGURES_Block F.dwg IVANNA.GOLDSBERRY 3/7/2016 5:35:24 PM

# APPENDIX D—UNDERGROUND STORAGE TANK RESIDUAL RISK TECHNICAL MEMORANDUM

**Technical Memorandum**
**Block F Underground Storage Tank**
**Lockheed Martin, Middle River Complex**
**Middle River, Maryland**
**May 2016**

*Block F UST Pipe Samples*

Detected concentrations in the Block F underground storage tank (UST) pipe samples (eight base samples and four sidewall samples plus one duplicate) were compared to Maryland Department of the Environment (MDE) Cleanup Standards for residential and non-residential soil and United States Environmental Protection Agency (USEPA) Regional Screening Levels (RSLs) for residential and industrial soil, and concentrations of benzo(a)pyrene equivalents (BaPEq) were compared to the remedial goal (RG) of 2,900 µg/kg. BaPEq were calculated for the carcinogenic polycyclic aromatic hydrocarbons (cPAHs) benzo(a)anthracene, benzo(a)pyrene, benzo(b)fluoranthene, benzo(k)fluoranthene, chrysene, dibenzo(a,h)anthracene, and indeno(1,2,3 cd)pyrene using a toxicity equivalence factor (TEF) approach based on the relative potency of each compound to that of benzo(a)pyrene (USEPA, 1993). The TEFs are used to convert each individual cPAH concentration into an equivalent concentration of benzo(a)pyrene; these individual concentrations are then summed. That summation is referred to as the BaPEq concentration. BaPEq concentrations were calculated two different ways. First, a BaPEq-half non-detect value was calculated using one-half of the detection limit of non-detected cPAHs to represent non-detected cPAH concentrations in the calculation. Second, the BaPEq-positive results value was calculated using a value of 0 for non-detected results. In either case, if all the cPAHs were not detected in a sample, then the sample quantitation limit for benzo(a)pyrene was used as the BaPEq concentration for that sample. The relevant data and calculations are presented in Table 1. The cPAHs were the only target analytes detected at concentrations exceeding the RSLs or MDE standards.

Concentrations of benzo(a)pyrene and BaPEq exceeded the MDE Cleanup Standards for residential and non-residential soil in one sample, F-VS-UST-P-3, and several additional detected concentrations of cPAHs exceeded the MDE Cleanup Standard for residential soil but did not exceed the MDE Cleanup Standard for non-residential soil (in samples F-VB-UST-P-8,

F-VS-UST-P-1-D, F-VS-UST-P-3, and F-VS-UST-P-4). Concentrations of BaPEq and benzo(a)pyrene exceeded the USEPA RSLs for industrial and residential soil in sample F-VS-UST-P-3, and concentrations of one or more cPAHs exceeded RSLs for residential soil only in samples F-VB-UST-P-8, F-VS-UST-P-1, F-VS-UST-P-1-D, F-VS-UST-P-2, F-VS-UST-P-3, and F-VS-UST-P-4. However, BaPEq concentrations reported for all samples collected are less than the RG for the industrial worker. It should be noted that RGs developed for BaPEq for construction workers or utility workers would exceed those developed for a typical industrial worker primarily because the exposure frequency/duration assumptions for the typical industrial worker are more conservative than those for the construction worker or utility worker. Thus, the BaPEq RG for the industrial worker is also protective of the construction worker or utility worker. An RG of 16,000 µg/kg was developed for construction workers exposed to BaPEq in subsurface soil in the 2012 human health risk assessment (HHRA) (Tetra Tech, 2012). A calculation showing how this RG was developed is attached. The construction worker RG is based on a target risk level of $1\times10^{-5}$.

Conservatively, a cancer risk estimate was developed assuming that an industrial worker might (hypothetically) be exposed to BaPEq concentrations in the soils. The analysis (presented below) demonstrates that the risk estimate would not exceed the MDE's cumulative cancer risk benchmark. This analysis is included for informational purposes only. Typical worker exposure to subsurface soils will be prevented by land use controls.

Because concentrations of individual cPAHs and the BaPEq (which represents the cPAHs) exceeded one or more screening criteria (i.e., the RSLs and MDE standards), the cPAHs were evaluated as chemicals of potential concern (COPC), and 95% upper confidence limits (UCLs) on the arithmetic mean for BaPEq (for both BaPEq-half non-detect and BaPEq-positive results) were calculated using USEPA ProUCL (Version 5.0.00) software (USEPA, 2013). Samples (all were subsurface soil collected greater than 2 feet below ground surface [bgs]) used for calculating the 95% UCLs and their BaPEq concentrations are listed in the following table:

| Sample | BaPEq-Half Non-Detect (µg/kg) | BaPEq-Positive Results (µg/kg) |
|---|---|---|
| F-VB-UST-P-1 | 0.75 U | 0.75 U |
| F-VB-UST-P-2 | 0.76 U | 0.76 U |
| F-VB-UST-P-3 | 0.78 U | 0.78 U |

| Sample | BaPEq-Half Non-Detect (µg/kg) | BaPEq-Positive Results (µg/kg) |
|---|---|---|
| F-VB-UST-P-4 | 0.77 U | 0.77 U |
| F-VB-UST-P-5 | 0.78 U | 0.78 U |
| F-VB-UST-P-6 | 0.75 U | 0.75 U |
| F-VB-UST-P-7 | 0.78 U | 0.78 U |
| F-VB-UST-P-8 | 150 | 150 |
| F-VS-UST-P-1-AVG | 20 | 19 |
| F-VS-UST-P-2 | 17 | 16 |
| F-VS-UST-P-3 | 1000 | 1000 |
| F-VS-UST-P-4 | 170 | 170 |

The 95% UCLs for the BaPEq are similar whether using one-half the value of non-detected results or positive results only in the calculation, and the higher of the UCLs, 810 µg/kg, was selected to represent the BaPEq. This UCL is less than the RG of 2,900 µg/kg. The ProUCL output file is included at the end of this appendix.

The 95% UCL was also used to calculate risk for an industrial worker, using the following equation:

$$\text{Cancer risk} = [(95\%\ \text{UCL} \times \text{Target risk}) / \text{Risk-based concentration}]$$

This simple risk estimation technique is presented in the User's Guide for the USEPA's RSL Tables. The target risk level in the calculation was set to $1\times10^{-6}$ (i.e., the RSL is set at a concentration representing the one-in-one million excess chance of developing cancer), and the RSL for industrial soil (USEPA, 2015) was used as the risk-based concentration.

For the industrial scenario, the residual risk would be:

$$\text{Cancer risk} = [(810\ \mu g/kg \times 1\times10^{-6}) / 290\ \mu g/kg] = 3\times10^{-6}.$$

The residual cancer risk for the industrial scenario ($3\times10^{-6}$) is less than MDE's cumulative cancer risk management benchmark of $1\times10^{-5}$.

*Block F UST Base Samples*

Detected concentrations in the Block F UST Area base samples (14 samples plus 1 duplicate) were compared to MDE and USEPA residential and industrial soil criteria. Detected results did not exceed the criteria, and cPAHs were not detected. Therefore, no COPCs were identified, and no further evaluation was necessary.

*References*

Tetra Tech, Inc. (Tetra Tech), 2012. Human Health Risk Assessment (HHRA) for Blocks D, E, F, G, and H Soils, Lockheed Martin Middle River Complex, 2323 Eastern Boulevard, Middle River, Maryland. September.

USEPA, 1993. Provisional Guidance for Quantitative Risk Assessment of Polycyclic Aromatic Hydrocarbons. EPA/600/R-93/089. Office of Research and Development, Washington, D.C. July.

USEPA, 2013. ProUCL Version 5.0.00 User Guide. Office of Research and Development, Washington, D.C. EPA/600/R 07/041. September.

USEPA, 2015. Regional Screening Levels for Chemical Contaminants at Superfund Sites. http://epa-prgs.USEPA.gov/chemicals/index.shtml. November.

**Site-specific**
**Construction Worker Equation Inputs for Soil - Unpaved Road Traffic**

| Variable | Value |
|---|---|
| TR (target cancer risk) unitless | 0.000001 |
| THQ (target hazard quotient) unitless | 1 |
| $EF_{cw}$ (exposure frequency - construction worker) day/yr | 250 |
| $ED_{cw}$ (exposure duration - construction worker) yr | 1 |
| $ET_{cw}$ (exposure time - construction worker) hr/day | 8 |
| LT (lifetime) yr | 70 |
| $BW_{cw}$ (body weight - construction worker) kg | 70 |
| $IR_{cw}$ (soil ingestion rate - construction worker) mg/day | 480 |
| $SA_{cw}$ (surface area - construction worker) $cm^2$/day | 3280 |
| $AF_{cw}$ (skin adherence factor - construction worker) $mg/cm^2$ | 0.3 |
| $AT_{cw}$ (averaging time - construction worker) | 365 |
| $EW_{cw}$ (overall duration of construction) weeks/year | 50 |
| $DW_{cw}$ (days worked - construction worker) days/week | 5 |
| $A_s$ ($PEF_{sc}$ - acres) | 0.5 |
| s (road surface silt content) % | 8.5 |
| $M_{dry}$ (road surface material moisture content under dry, uncontrolled conditions) % | 0.2 |
| p (days per year with at least .01" of precipitation) days/year | 140 |
| $L_R$ (length of road segment) ft | 147.58077 |
| $W_R$ (width of road segment) ft | 20 |
| number of cars | 20 |
| number of trucks | 10 |
| tons/car | 2 |
| tons/truck | 20 |
| $F_D$ Unitless Dispersion Correction Factor | 0.185837208 |
| $t_c$ (overall duration of construction) hours | 8400 |
| distance (road length) km/day | 0.04498 |
| $T_t$ (overall duration of traffic) s | 7200000 |
| total number of vehicles | 30 |
| $A_R$ (surface area of contaminated road segment) $m^2$ | 274.21393 |
| W (mean vehicle weight) tons | 8 |
| ΣVKT (sum of fleet vehicle km traveled) km | 337.35 |
| $Q/C_{sr}$ (inverse of the ratio of the 1-h. geometric mean air concentration to the emissio | 23.01785 |
| $PEF_{sc}$ (particulate emission factor) $m^3/kg$ | 1428031.096 |
| A (Dispersion Constant) | 12.9351 |
| B (Dispersion Constant) | 5.7383 |
| C (Dispersion Constant) | 71.7711 |
| $A_{surf}$ (areal extent of site) $m^2$ | |
| $A_s$ ($VF_{ulim-sc}$ acres) | 0.5 |
| T (temperature) °C | 25 |
| foc (fraction organic carbon in soil) g/g | 0.006 |
| $\rho_b$ (dry soil bulk density) $g/cm^3$ | 1.5 |
| $\rho_s$ (soil particle density) $g/cm^3$ | 2.65 |
| $\theta_w$ (water-filled soil porosity) $L_{water}/L_{soil}$ | 0.15 |
| A (VF Dispersion Constant) | 2.4538 |
| B (VF Dispersion Constant) | 17.566 |
| C (VF Dispersion Constant) | 189.0426 |
| $Q/C_{sa}$ (inverse of the ratio of the geometric mean air concentration to the emission flu | 14.31407 |
| n (total soil porosity) $L_{pore}/L_{soil}$ | 0.43396 |
| $\theta_a$ (air-filled soil porosity) $L_{air}/L_{soil}$ | 0.28396 |
| $A_s$ ($VF_{mlim-sc}$ acres) | 0.5 |
| $\rho_b$ (dry soil bulk density) $g/cm^3$ | 1.5 |
| $d_s$ (average source depth) m | . |
| $Q/C_{vol}$ (inverse of the ratio of the geometric mean air concentration to the emission fl | 14.31407 |
| $VF_{mlim-sc}$ (volitization factor) $m^3_{air}/kg_{soil}$ | . |

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**Site-specific**
**Construction Worker Screening Levels (RSL) for Soil - Unpaved Road Traffic**
ca=Cancer, nc=Noncancer, ca* (Where nc SL < 100 x ca SL),
ca** (Where nc SL < 10 x ca SL), max=SL exceeds ceiling limit (see User's Guide), sat=SL exceeds csat,
Smax=Soil SL exceeds ceiling limit and has been substituted with the max value (see User's Guide),
Ssat=Soil inhalation SL exceeds csat and has been substituted with the csat

| Chemical | CAS Number | Mutagen? | VOC? | Ingestion SF (mg/kg-day)$^{-1}$ | SFO Ref | Inhalation Unit Risk (ug/m$^3$)$^{-1}$ | IUR Ref | Subchronic RfD (mg/kg-day) | Subchronic RfD Ref | Subchronic RfC (mg/m$^3$) | Subchronic RfC Ref | GIABS | ABS | RBA | Volatilization Factor (m$^3$/kg) | Soil Saturation Concentration (mg/kg) | Particulate Emission Factor (m$^3$/kg) | Ingestion SL TR=1.0E-6 (mg/kg) | Dermal SL TR=1.0E-6 (mg/kg) | Inhalation SL TR=1.0E-6 (mg/kg) | Carcinogenic SL TR=1.0E-6 (mg/kg) | Ingestion SL THQ=1 (mg/kg) | Dermal SL THQ=1 (mg/kg) | Inhalation SL THQ=1 (mg/kg) | Noncarcinogenic SL THI=1 (mg/kg) | Screening Level (mg/kg) |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| Benzo[a]pyrene | 50-32-8 | Yes | No | 7.30E+00 | I | 1.10E-03 | C | - | | - | | 1 | 0.13 | 1 | - | - | 1.43E+06 | 2.04E+00 | 7.66E+00 | 3.98E+02 | 1.61E+00 | - | - | - | - | 1.61E+00 ca |

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# APPENDIX E—ANALYTICAL LABORATORY REPORT

# LABORATORY REPORTS

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

# ANALYTICAL REPORT

Job Number: 240-57597-1

Job Description: MRC Soil Remedy Block F UST

For:
Tetra Tech, Inc.
20251 Century Blvd
Suite 200
Germantown, MD 20874

Attention: Tony Apanavage



Approved for release.
John McFadden
Project Manager I
11/19/2015 10:29 AM

John McFadden, Project Manager I
4101 Shuffel Street NW, North Canton, OH, 44720
john.mcfadden@testamericainc.com
11/19/2015
Revision: 1

cc: Samantha Brenner
Kelly Carper
Tobrena Sedlmyer

The test results in this report meet all NELAP requirements for parameters for which accreditation is required or available. Any exceptions to the NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory. This report is confidential and is intended for the sole use of TestAmerica and its client. All questions regarding this report should be directed to the TestAmerica Project Manager who has signed this report.

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# CASE NARRATIVE

## Client: Tetra Tech, Inc.

## Project: MRC Soil Remedy Block F UST

## Report Number: 240-57597-1

With the exceptions noted as flags or footnotes, standard analytical protocols were followed in the analysis of the samples and no problems were encountered or anomalies observed. In addition all laboratory quality control samples were within established control limits, with any exceptions noted below. Each sample was analyzed to achieve the lowest possible reporting limit within the constraints of the method. In some cases, due to interference or analytes present at high concentrations, samples were diluted. For diluted samples, the reporting limits are adjusted relative to the dilution required.

TestAmerica Canton attests to the validity of the laboratory data generated by TestAmerica facilities reported herein. All analyses performed by TestAmerica facilities were done using established laboratory SOPs that incorporate QA/QC procedures described in the application methods. TestAmerica's operations groups have reviewed the data for compliance with the laboratory QA/QC plan, and data have been found to be compliant with laboratory protocols unless otherwise noted below.

The test results in this report meet all NELAP requirements for parameters for which accreditation is required or available. Any exceptions to NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory.

Calculations are performed before rounding to avoid round-off errors in calculated results.

All holding times were met and proper preservation noted for the methods performed on these samples, unless otherwise detailed in the individual sections below.

All solid sample results are reported on an "as received" basis unless otherwise indicated by the presence of a % solids value in the method header.

This laboratory report is confidential and is intended for the sole use of TestAmerica and its client.

**RECEIPT**

The samples were received on 11/6/2015 10:20 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperatures of the 2 coolers at receipt time were 0.9º C and 2.2º C.

**VOLATILE ORGANIC COMPOUNDS (GCMS)**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for volatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8260C. The samples were prepared and analyzed on 11/10/2015 and 11/11/2015.

2-Butanone (MEK), Naphthalene and Toluene were detected in method blank MB 240-206267/6 at levels that were above the method detection limit but below the reporting limit. The values should be considered estimates, and have been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

2-Butanone (MEK) was detected in method blank MB 240-206089/1-A at a level that was above the method detection limit but below the reporting limit. The value should be considered an estimate, and has been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

The continuing calibration verification (CCV) associated with batch 240-206051 recovered above the upper control limit for 1,1,1-Trichloroethane; 1,1,2-Trichloro-1,2,2-trifluoroethane; 2,2-Dichloropropane; Carbon tetrachloride; Dichloro-difluoromethane and Trichlorofluoromethane. The samples associated with this CCV were non-detects for the affected analytes; therefore, the data have been reported.

The continuing calibration verification (CCV) analyzed in batch 240-206267 was outside the method criteria for the following analyte(s): 2-Hexanone, Acetone, 2-Methyl-2-propanol, 2-Chloroethyl Vinyl Ether, 4-Methyl-2-Pentanone, 2,2-Dichloropropane, and 2-Butanone. An MRL standard at or below the reporting limit (RL) was analyzed with the affected samplesF-VB-UST-DUP (240-57597-15), (CCV 240-206267/4), (CCVIS 240-206267/3), (LCS 240-206267/5), (MB 240-206267/6), (240-57597-B-15-C MS) and (240-57597-B-15-D MSD) and found to be acceptable. As indicated in the reference method, sample analysis may proceed; however, any detection for the affected analyte(s) is considered estimated.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**VOLATILE ORGANIC COMPOUNDS (GCMS)**

Sample TB-110515 (240-57597-16) was analyzed for volatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8260C. The samples were analyzed on 11/12/2015.

2-Butanone (MEK) was detected in method blank MB 240-206461/6 at a level that was above the method detection limit but below the reporting limit. The value should be considered an estimate, and has been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

The continuing calibration verification (CCV) for analytical batch 206461 exceeded control criteria for Hexachlorobutadiene. The samples associated with this CCV were non-detect for the affected analyte. In accordance with the laboratory SOP, a low level CCV at the reporting limit (labeled as an MRL) was analyzed and the affected compound was detected; therefore the data has been reported. No further corrective action was required. TB-110515 (240-57597-16)

2-Chloroethyl vinyl ether cannot be reliably recovered in an acid preserved sample. TB-110515 (240-57597-16)

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**SEMIVOLATILE ORGANIC COMPOUNDS (GCMS)**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for semivolatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8270D. The samples were prepared on 11/09/2015 and analyzed on 11/11/2015.

The analyte benzaldehyde had an Rf below method recommendations in analytical batch 206233. An LODV was analyzed to demonstrate sensitivity for this analyte at the reporting limit for the following samples:F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14), F-VB-UST-DUP (240-57597-15), (LCS 240-205866/9-A), (MB 240-205866/8-A), (240-57597-D-15-E MS) and (240-57597-D-15-F MSD).

Two analytes had Rfs below method recommendations: Benzaldehyde and Hexachlorocyclopentadiene. A LODV was analyzed to demonstrate sensitivity for these analytes at the reporting limit for the following samples: F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9) and F-VB-UST-10-14 (240-57597-10).

The continuing calibration verification (CCV) associated with analytical batch 206233 recovered above the upper control limit for N-Nitrosodi-n-propylamine. The samples associated with this CCV were non-detect for this analyte; therefore, the data have been reported. The following samples are impacted: F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14), F-VB-UST-DUP (240-57597-15), (MB 240-205866/8-A), (240-57597-D-15-E MS) and (240-57597-D-15-F MSD).

The continuing calibration verification (CCV) for analytical batch 206233 was below method recommendations for the analyte benzaldehyde (>20%D). Benzaldehyde was present in sample F-VB-UST-11-14 (240-57597-11) at an estimated amount of 19ug/kg J. Had the response of benzaldehyde been within method recommendations the amount of benzaldehyde measured in the sample would likely have still been a small estimated concentration (J flag). Therefore, no additional analysis or corrective action was initiated. The results have been flagged accordingly and reported.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**GASOLINE RANGE ORGANICS**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for gasoline range organics in accordance with EPA SW-846 Method 8015C - GRO. The samples were prepared and analyzed on 11/11/2015 and 11/12/2015.

The matrix spike duplicate (MSD) precision for 206211 was outside control limits. Sample matrix interference and/or non-homogeneity are suspected because the associated laboratory control sample (LCS) precision was within acceptance limits.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**DIESEL RANGE ORGANICS (DRO)**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12),

F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for diesel range organics (DRO) in accordance with EPA SW-846 Method 8015C - DRO. The samples were prepared on 11/09/2015 and analyzed on 11/10/2015 and 11/11/2015.

Total Petroleum Hydrocarbons (C10-C32) was detected in method blank MB 240-205788/22-A at a level exceeding the reporting limit. If the associated sample reported a result above the MDL and/or RL, the result has been flagged. .

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**POLYCHLORINATED BIPHENYLS (PCBS)**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for polychlorinated biphenyls (PCBs) in accordance with EPA SW-846 Method 8082A. The samples were prepared on 11/09/2015 and analyzed on 11/11/2015.

Surrogates are added during the extraction process prior to dilution. When the sample dilution is 5X or greater, surrogate recoveries are diluted out and no corrective action is required. All of the samples in this data set analyzed for PCBs were subjected to the sulfuric acid cleanup procedure before instrumental analysis, per EPA Method 3665A.

The following samples required a tetrabutylammonium sulfite (TBA) clean-up to reduce matrix interferences caused by sulfur: F-VB-UST-5-14 (240-57597-5) and F-VB-UST-13-14 (240-57597-13).Reagents 2266653, 1848564 and 2283702.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**PERCENT SOLIDS**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for percent solids in accordance with EPA Method 160.3 MOD. The samples were analyzed on 11/09/2015.

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 200091

Lab Sample ID: STD5 240-200091/2 IC Client Sample ID:

Date Analyzed: 10/02/15 10:48 Lab File ID: 51002002.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.92 | Poor chromatography | gruberj | 10/02/15 11:35 |
| N-Nitrosodimethylamine | 3.41 | Poor chromatography | gruberj | 10/02/15 11:35 |
| Pyridine | 3.47 | Poor chromatography | gruberj | 10/02/15 11:35 |

Lab Sample ID: STD4 240-200091/3 IC Client Sample ID:

Date Analyzed: 10/02/15 11:12 Lab File ID: 51002003.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.93 | Poor chromatography | gruberj | 10/02/15 11:37 |
| N-Nitrosodimethylamine | 3.40 | Poor chromatography | gruberj | 10/02/15 11:37 |
| Pyridine | 3.47 | Poor chromatography | gruberj | 10/02/15 11:37 |
| Benzoic acid | 7.04 | Poor chromatography | gruberj | 10/02/15 11:37 |

Lab Sample ID: STD3 240-200091/4 IC Client Sample ID:

Date Analyzed: 10/02/15 11:37 Lab File ID: 51002004.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.93 | Poor chromatography | gruberj | 10/02/15 12:48 |
| N-Nitrosodimethylamine | 3.41 | Poor chromatography | gruberj | 10/02/15 12:48 |
| Pyridine | 3.47 | Poor chromatography | gruberj | 10/02/15 12:48 |
| Benzoic acid | 7.03 | Poor chromatography | gruberj | 10/02/15 12:48 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 200091

Lab Sample ID: STD2 240-200091/5 IC Client Sample ID:

Date Analyzed: 10/02/15 12:01 Lab File ID: 51002005.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.96 | Poor chromatography | gruberj | 10/02/15 12:50 |
| N-Nitrosodimethylamine | 3.38 | Poor chromatography | gruberj | 10/02/15 12:50 |
| Pyridine | 3.47 | Poor chromatography | gruberj | 10/02/15 12:50 |

Lab Sample ID: STD9 240-200091/7 IC Client Sample ID:

Date Analyzed: 10/02/15 12:50 Lab File ID: 51002007.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.93 | Poor chromatography | gruberj | 10/02/15 14:00 |
| N-Nitrosodimethylamine | 3.41 | Poor chromatography | gruberj | 10/02/15 14:00 |
| Pyridine | 3.47 | Poor chromatography | gruberj | 10/02/15 14:00 |
| Benzoic acid | 7.13 | Poor chromatography | gruberj | 10/02/15 14:00 |
| Caprolactam | 7.71 | Poor chromatography | gruberj | 10/02/15 14:00 |

Lab Sample ID: STD8 240-200091/8 IC Client Sample ID:

Date Analyzed: 10/02/15 13:15 Lab File ID: 51002008.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.96 | Poor chromatography | gruberj | 10/02/15 14:02 |
| N-Nitrosodimethylamine | 3.41 | Poor chromatography | gruberj | 10/02/15 14:02 |
| Pyridine | 3.47 | Poor chromatography | gruberj | 10/02/15 14:02 |
| Benzoic acid | 7.11 | Poor chromatography | gruberj | 10/02/15 14:02 |
| Caprolactam | 7.70 | Poor chromatography | gruberj | 10/02/15 14:02 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 200091

Lab Sample ID: STD7 240-200091/9 IC Client Sample ID:

Date Analyzed: 10/02/15 13:39 Lab File ID: 51002009.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.93 | Poor chromatography | gruberj | 10/02/15 14:04 |
| N-Nitrosodimethylamine | 3.41 | Poor chromatography | gruberj | 10/02/15 14:04 |
| Pyridine | 3.47 | Poor chromatography | gruberj | 10/02/15 14:04 |
| Benzoic acid | 7.10 | Poor chromatography | gruberj | 10/02/15 14:04 |
| Caprolactam | 7.69 | Poor chromatography | gruberj | 10/02/15 14:04 |

Lab Sample ID: STD6 240-200091/10 ICIS Client Sample ID:

Date Analyzed: 10/02/15 14:04 Lab File ID: 51002010.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.93 | Poor chromatography | gruberj | 10/02/15 14:27 |
| N-Nitrosodimethylamine | 3.41 | Poor chromatography | gruberj | 10/02/15 14:27 |
| Pyridine | 3.47 | Poor chromatography | gruberj | 10/02/15 14:27 |
| Caprolactam | 7.68 | Poor chromatography | gruberj | 10/02/15 14:27 |

Lab Sample ID: ICV 240-200091/11 Client Sample ID:

Date Analyzed: 10/02/15 14:28 Lab File ID: 51002011.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.96 | Poor chromatography | gruberj | 10/02/15 15:03 |
| N-Nitrosodimethylamine | 3.41 | Poor chromatography | gruberj | 10/02/15 15:03 |
| Pyridine | 3.47 | Poor chromatography | gruberj | 10/02/15 15:03 |
| Caprolactam | 7.68 | Poor chromatography | gruberj | 10/02/15 15:03 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 206256

Lab Sample ID: CCV 240-206256/2 CCVIS Client Sample ID:

Date Analyzed: 11/11/15 09:34 Lab File ID: 51111002.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.75 | Poor chromatography | gruberj | 11/11/15 09:57 |
| N-Nitrosodimethylamine | 3.28 | Poor chromatography | gruberj | 11/11/15 09:57 |
| Pyridine | 3.31 | Poor chromatography | gruberj | 11/11/15 09:57 |

Lab Sample ID: 240-57597-7 Client Sample ID: F-VB-UST-7-14

Date Analyzed: 11/11/15 13:44 Lab File ID: 51111012.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 2,4,6-Tribromophenol (Surr) | 9.45 | Poor chromatography | gruberj | 11/11/15 14:04 |

Lab Sample ID: 240-57597-9 Client Sample ID: F-VB-UST-9-14

Date Analyzed: 11/11/15 14:33 Lab File ID: 51111014.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Phenol-d5 (Surr) | 5.83 | Poor chromatography | gruberj | 11/12/15 12:45 |
| 2,4,6-Tribromophenol (Surr) | 9.45 | Poor chromatography | gruberj | 11/12/15 12:45 |

Lab Sample ID: 240-57597-10 Client Sample ID: F-VB-UST-10-14

Date Analyzed: 11/11/15 14:57 Lab File ID: 51111015.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Phenol-d5 (Surr) | 5.83 | Poor chromatography | gruberj | 11/12/15 12:46 |
| 2,4,6-Tribromophenol (Surr) | 9.45 | Poor chromatography | gruberj | 11/12/15 12:46 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP9 Analysis Batch Number: 204569

Lab Sample ID: STD5 240-204569/2 IC Client Sample ID:

Date Analyzed: 10/30/15 13:19 Lab File ID: 51030002.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.05 | Poor chromatography | ulmanm | 10/30/15 14:26 |
| N-Nitrosodimethylamine | 3.56 | Poor chromatography | ulmanm | 10/30/15 14:26 |
| Pyridine | 3.62 | Poor chromatography | ulmanm | 10/30/15 14:26 |
| Benzoic acid | 7.24 | Poor chromatography | ulmanm | 10/30/15 14:26 |
| 4-Chloroaniline | 7.57 | Poor chromatography | ulmanm | 10/30/15 14:26 |
| Di-n-octyl phthalate | 13.17 | Poor chromatography | ulmanm | 10/30/15 14:26 |

Lab Sample ID: STD4 240-204569/3 IC Client Sample ID:

Date Analyzed: 10/30/15 13:44 Lab File ID: 51030003.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.05 | Poor chromatography | ulmanm | 10/30/15 14:30 |
| N-Nitrosodimethylamine | 3.56 | Poor chromatography | ulmanm | 10/30/15 14:30 |
| Pyridine | 3.62 | Poor chromatography | ulmanm | 10/30/15 14:30 |
| Benzoic acid | 7.25 | Poor chromatography | ulmanm | 10/30/15 14:30 |
| Di-n-octyl phthalate | 13.16 | Poor chromatography | ulmanm | 10/30/15 14:30 |
| Indeno[1,2,3-cd]pyrene | 16.42 | Poor chromatography | ulmanm | 10/30/15 14:30 |

Lab Sample ID: STD3 240-204569/4 IC Client Sample ID:

Date Analyzed: 10/30/15 14:10 Lab File ID: 51030004.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.04 | Poor chromatography | ulmanm | 10/30/15 14:43 |
| Pyridine | 3.62 | Poor chromatography | ulmanm | 10/30/15 14:43 |
| Benzoic acid | 7.23 | Poor chromatography | ulmanm | 10/30/15 14:43 |
| Di-n-octyl phthalate | 13.16 | Poor chromatography | ulmanm | 10/30/15 14:43 |
| Indeno[1,2,3-cd]pyrene | 16.41 | Poor chromatography | ulmanm | 10/30/15 14:43 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP9 Analysis Batch Number: 204569

Lab Sample ID: STD2 240-204569/5 IC Client Sample ID:

Date Analyzed: 10/30/15 14:35 Lab File ID: 51030005.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.06 | Poor chromatography | ulmanm | 10/30/15 15:14 |
| N-Nitrosodimethylamine | 3.56 | Poor chromatography | ulmanm | 10/30/15 15:14 |
| Pyridine | 3.62 | Poor chromatography | ulmanm | 10/30/15 15:14 |
| 4-Nitroaniline | 9.49 | Poor chromatography | ulmanm | 10/30/15 15:14 |
| Di-n-octyl phthalate | 13.16 | Poor chromatography | ulmanm | 10/30/15 15:14 |
| Dibenz(a,h)anthracene | 16.42 | Poor chromatography | ulmanm | 10/30/15 15:14 |
| Indeno[1,2,3-cd]pyrene | 16.42 | Poor chromatography | ulmanm | 10/30/15 15:14 |
| Benzo[g,h,i]perylene | 17.04 | Poor chromatography | ulmanm | 10/30/15 15:14 |

Lab Sample ID: STD1 240-204569/6 IC Client Sample ID:

Date Analyzed: 10/30/15 15:01 Lab File ID: 51030006.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Benzo[b]fluoranthene | 13.85 | Poor chromatography | ulmanm | 10/30/15 15:45 |
| Benzo[a]pyrene | 14.34 | Poor chromatography | ulmanm | 10/30/15 15:45 |
| Indeno[1,2,3-cd]pyrene | 16.41 | Poor chromatography | ulmanm | 10/30/15 15:45 |
| Dibenz(a,h)anthracene | 16.42 | Poor chromatography | ulmanm | 10/30/15 15:45 |
| Benzo[g,h,i]perylene | 17.03 | Poor chromatography | ulmanm | 10/30/15 15:45 |

Lab Sample ID: STD9 240-204569/7 IC Client Sample ID:

Date Analyzed: 10/30/15 15:27 Lab File ID: 51030007.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.05 | Poor chromatography | ulmanm | 10/30/15 16:05 |
| Pyridine | 3.62 | Poor chromatography | ulmanm | 10/30/15 16:05 |
| Benzoic acid | 7.29 | Poor chromatography | ulmanm | 10/30/15 16:05 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP9 Analysis Batch Number: 204569

Lab Sample ID: STD8 240-204569/8 IC Client Sample ID:

Date Analyzed: 10/30/15 15:52 Lab File ID: 51030008.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.05 | Poor chromatography | ulmanm | 10/30/15 16:32 |
| Caprolactam | 7.88 | Poor chromatography | ulmanm | 10/30/15 16:32 |

Lab Sample ID: STD7 240-204569/9 IC Client Sample ID:

Date Analyzed: 10/30/15 16:17 Lab File ID: 51030009.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.05 | Poor chromatography | ulmanm | 10/31/15 07:11 |

Lab Sample ID: STD6 240-204569/10 ICIS Client Sample ID:

Date Analyzed: 10/30/15 16:42 Lab File ID: 51030010.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.06 | Poor chromatography | ulmanm | 10/31/15 07:17 |
| Pyridine | 3.62 | Poor chromatography | ulmanm | 10/31/15 07:17 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP9 Analysis Batch Number: 206233

Lab Sample ID: CCV 240-206233/2 CCVIS Client Sample ID:

Date Analyzed: 11/11/15 07:49 Lab File ID: 51111002.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Benzoic acid | 7.15 | Poor chromatography | ulmanm | 11/11/15 08:27 |

Lab Sample ID: 240-57597-12 Client Sample ID: F-VB-UST-12-14

Date Analyzed: 11/11/15 15:14 Lab File ID: 51111020.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 2,4,6-Tribromophenol (Surr) | 9.59 | Poor chromatography | ulmanm | 11/12/15 08:11 |

Lab Sample ID: 240-57597-13 Client Sample ID: F-VB-UST-13-14

Date Analyzed: 11/11/15 15:39 Lab File ID: 51111021.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 2,4,6-Tribromophenol (Surr) | 9.59 | Poor chromatography | ulmanm | 11/12/15 08:13 |

Lab Sample ID: 240-57597-14 Client Sample ID: F-VB-UST-14-14

Date Analyzed: 11/11/15 16:03 Lab File ID: 51111022.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Phenol-d5 (Surr) | 5.96 | Poor chromatography | ulmanm | 11/12/15 08:15 |
| 2,4,6-Tribromophenol (Surr) | 9.59 | Poor chromatography | ulmanm | 11/12/15 08:15 |

Lab Sample ID: 240-57597-15 Client Sample ID: F-VB-UST-DUP

Date Analyzed: 11/11/15 16:28 Lab File ID: 51111023.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Phenol-d5 (Surr) | 5.97 | Poor chromatography | ulmanm | 11/12/15 08:26 |
| 2,4,6-Tribromophenol (Surr) | 9.59 | Poor chromatography | ulmanm | 11/12/15 08:26 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP9 Analysis Batch Number: 206233

Lab Sample ID: 240-57597-15 MS Client Sample ID: F-VB-UST-DUP MS

Date Analyzed: 11/11/15 16:52 Lab File ID: 51111024.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Phenol-d5 (Surr) | 5.96 | Poor chromatography | ulmanm | 11/12/15 08:31 |
| 3 & 4 Methylphenol | 6.64 | Poor chromatography | ulmanm | 11/12/15 08:31 |
| 4-Chloro-3-methylphenol | 7.89 | Poor chromatography | ulmanm | 11/12/15 08:31 |
| 2,4,5-Trichlorophenol | 8.33 | Poor chromatography | ulmanm | 11/12/15 08:31 |
| 4-Nitrophenol | 8.98 | Poor chromatography | ulmanm | 11/12/15 08:31 |
| 2,4,6-Tribromophenol (Surr) | 9.59 | Poor chromatography | ulmanm | 11/12/15 08:31 |
| Pentachlorophenol | 10.02 | Poor chromatography | ulmanm | 11/12/15 08:31 |

Lab Sample ID: 240-57597-15 MSD Client Sample ID: F-VB-UST-DUP MSD

Date Analyzed: 11/11/15 17:17 Lab File ID: 51111025.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 2-Methylphenol | 6.50 | Poor chromatography | ulmanm | 11/12/15 08:35 |
| 3 & 4 Methylphenol | 6.63 | Poor chromatography | ulmanm | 11/12/15 08:35 |
| 4-Chloro-3-methylphenol | 7.87 | Poor chromatography | ulmanm | 11/12/15 08:35 |
| 2,4,5-Trichlorophenol | 8.30 | Poor chromatography | ulmanm | 11/12/15 08:35 |
| 2,4,6-Tribromophenol (Surr) | 9.59 | Poor chromatography | ulmanm | 11/12/15 08:35 |

GC VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: AFID Analysis Batch Number: 194508

Lab Sample ID: STD1 240-194508/7 IC Client Sample ID:

Date Analyzed: 08/24/15 13:59 Lab File ID: A0082107.D GC Column: RTX-VRX ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Trifluorotoluene (Surr) | 10.34 | Baseline Smoothing | bosworthh | 08/25/15 10:02 |
| Gasoline Range Organics [C6 - C10] | 11.46 | Baseline Smoothing | bosworthh | 08/25/15 10:02 |
| C6-C12 | 12.73 | Baseline Smoothing | bosworthh | 08/25/15 10:02 |

DIESEL RANGE ORGANICS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP5R Analysis Batch Number: 206105

Lab Sample ID: CCV 240-206105/7 CCVRT Client Sample ID:

Date Analyzed: 11/10/15 16:28 Lab File ID: P5B11007.D GC Column: ID:

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| o-Terphenyl (Surr) | 9.83 | Other | bolgrind | 11/10/15 17:01 |

DIESEL RANGE ORGANICS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP5R Analysis Batch Number: 206356

Lab Sample ID: 240-57597-15 MSD Client Sample ID: F-VB-UST-DUP MSD

Date Analyzed: 11/11/15 21:39 Lab File ID: P5B11114.D GC Column: ID:

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| o-Terphenyl (Surr) | 9.83 | Other | bolgrind | 11/12/15 13:52 |

# SAMPLE SUMMARY

Client: Tetra Tech, Inc. Job Number: 240-57597-1

| Lab Sample ID | Client Sample ID | Client Matrix | Date/Time Sampled | Date/Time Received |
|---|---|---|---|---|
| 240-57597-1 | F-VB-UST-1-14 | Solid | 11/05/2015 0900 | 11/06/2015 1020 |
| 240-57597-2 | F-VB-UST-2-14 | Solid | 11/05/2015 0902 | 11/06/2015 1020 |
| 240-57597-3 | F-VB-UST-3-14 | Solid | 11/05/2015 0904 | 11/06/2015 1020 |
| 240-57597-4 | F-VB-UST-4-14 | Solid | 11/05/2015 0906 | 11/06/2015 1020 |
| 240-57597-5 | F-VB-UST-5-14 | Solid | 11/05/2015 0908 | 11/06/2015 1020 |
| 240-57597-6 | F-VB-UST-6-14 | Solid | 11/05/2015 0910 | 11/06/2015 1020 |
| 240-57597-7 | F-VB-UST-7-14 | Solid | 11/05/2015 0912 | 11/06/2015 1020 |
| 240-57597-8 | F-VB-UST-8-14 | Solid | 11/05/2015 0914 | 11/06/2015 1020 |
| 240-57597-9 | F-VB-UST-9-14 | Solid | 11/05/2015 0916 | 11/06/2015 1020 |
| 240-57597-10 | F-VB-UST-10-14 | Solid | 11/05/2015 0918 | 11/06/2015 1020 |
| 240-57597-11 | F-VB-UST-11-14 | Solid | 11/05/2015 0920 | 11/06/2015 1020 |
| 240-57597-12 | F-VB-UST-12-14 | Solid | 11/05/2015 0922 | 11/06/2015 1020 |
| 240-57597-13 | F-VB-UST-13-14 | Solid | 11/05/2015 0924 | 11/06/2015 1020 |
| 240-57597-14 | F-VB-UST-14-14 | Solid | 11/05/2015 0926 | 11/06/2015 1020 |
| 240-57597-15 | F-VB-UST-DUP | Solid | 11/05/2015 0000 | 11/06/2015 1020 |
| 240-57597-16 | TB-110515 | Water | 11/05/2015 0000 | 11/06/2015 1020 |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-57597-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-57597-1** | **F-VB-UST-1-14** | | | | | |
| 2-Butanone (MEK) | | 3.9 | J B | 20 | ug/Kg | 8260C |
| Acetophenone | | 16 | J | 120 | ug/Kg | 8270D |
| Bis(2-ethylhexyl) phthalate | | 31 | J | 83 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 16 | J B | 19 | mg/Kg | 8015C |
| Percent Solids | | 84 | | 0.10 | % | Moisture |
| Percent Moisture | | 16 | | 0.10 | % | Moisture |
| **240-57597-2** | **F-VB-UST-2-14** | | | | | |
| 2-Butanone (MEK) | | 4.0 | J B | 22 | ug/Kg | 8260C |
| Total Petroleum Hydrocarbons (C10-C32) | | 10 | J B | 20 | mg/Kg | 8015C |
| Percent Solids | | 85 | | 0.10 | % | Moisture |
| Percent Moisture | | 15 | | 0.10 | % | Moisture |
| **240-57597-3** | **F-VB-UST-3-14** | | | | | |
| 2-Butanone (MEK) | | 3.9 | J B | 23 | ug/Kg | 8260C |
| Percent Solids | | 82 | | 0.10 | % | Moisture |
| Percent Moisture | | 18 | | 0.10 | % | Moisture |
| **240-57597-4** | **F-VB-UST-4-14** | | | | | |
| 2-Butanone (MEK) | | 4.0 | J B | 22 | ug/Kg | 8260C |
| Acetophenone | | 13 | J | 120 | ug/Kg | 8270D |
| Bis(2-ethylhexyl) phthalate | | 33 | J | 83 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 7.6 | J B | 20 | mg/Kg | 8015C |
| Percent Solids | | 84 | | 0.10 | % | Moisture |
| Percent Moisture | | 16 | | 0.10 | % | Moisture |
| **240-57597-5** | **F-VB-UST-5-14** | | | | | |
| 2-Butanone (MEK) | | 4.3 | J B | 23 | ug/Kg | 8260C |
| Acetone | | 13 | J | 23 | ug/Kg | 8260C |
| Naphthalene | | 0.41 | J | 5.8 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 38 | J | 82 | ug/Kg | 8270D |
| Gasoline Range Organics [C6 - C10] | | 77 | J | 120 | ug/Kg | 8015C |
| Percent Solids | | 85 | | 0.10 | % | Moisture |
| Percent Moisture | | 15 | | 0.10 | % | Moisture |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-57597-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-57597-6** | **F-VB-UST-6-14** | | | | | |
| 2-Butanone (MEK) | | 3.8 | J B | 23 | ug/Kg | 8260C |
| Naphthalene | | 0.45 | J | 5.8 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 32 | J | 83 | ug/Kg | 8270D |
| Percent Solids | | 83 | | 0.10 | % | Moisture |
| Percent Moisture | | 17 | | 0.10 | % | Moisture |
| **240-57597-7** | **F-VB-UST-7-14** | | | | | |
| 2-Butanone (MEK) | | 4.1 | J B | 22 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 32 | J | 85 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 16 | J B | 20 | mg/Kg | 8015C |
| Percent Solids | | 84 | | 0.10 | % | Moisture |
| Percent Moisture | | 16 | | 0.10 | % | Moisture |
| **240-57597-8** | **F-VB-UST-8-14** | | | | | |
| 2-Butanone (MEK) | | 4.3 | J B | 23 | ug/Kg | 8260C |
| Toluene | | 0.33 | J | 5.8 | ug/Kg | 8260C |
| Percent Solids | | 82 | | 0.10 | % | Moisture |
| Percent Moisture | | 18 | | 0.10 | % | Moisture |
| **240-57597-9** | **F-VB-UST-9-14** | | | | | |
| 2-Butanone (MEK) | | 4.4 | J B | 23 | ug/Kg | 8260C |
| Acetone | | 12 | J | 23 | ug/Kg | 8260C |
| Toluene | | 0.36 | J | 5.8 | ug/Kg | 8260C |
| Acetophenone | | 13 | J | 120 | ug/Kg | 8270D |
| Bis(2-ethylhexyl) phthalate | | 40 | J | 87 | ug/Kg | 8270D |
| Percent Solids | | 81 | | 0.10 | % | Moisture |
| Percent Moisture | | 19 | | 0.10 | % | Moisture |
| **240-57597-10** | **F-VB-UST-10-14** | | | | | |
| 2-Butanone (MEK) | | 3.5 | J B | 22 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 29 | J | 81 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 13 | J B | 19 | mg/Kg | 8015C |
| Percent Solids | | 87 | | 0.10 | % | Moisture |
| Percent Moisture | | 13 | | 0.10 | % | Moisture |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-57597-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-57597-11** | **F-VB-UST-11-14** | | | | | |
| 2-Butanone (MEK) | | 3.9 | J B | 24 | ug/Kg | 8260C |
| Acetophenone | | 38 | J | 120 | ug/Kg | 8270D |
| Benzaldehyde | | 19 | J | 120 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 9.0 | J B | 20 | mg/Kg | 8015C |
| Percent Solids | | 83 | | 0.10 | % | Moisture |
| Percent Moisture | | 17 | | 0.10 | % | Moisture |
| **240-57597-12** | **F-VB-UST-12-14** | | | | | |
| 2-Butanone (MEK) | | 3.7 | J B | 22 | ug/Kg | 8260C |
| Toluene | | 0.30 | J | 5.5 | ug/Kg | 8260C |
| Total Petroleum Hydrocarbons (C10-C32) | | 30 | B | 20 | mg/Kg | 8015C |
| Percent Solids | | 84 | | 0.10 | % | Moisture |
| Percent Moisture | | 16 | | 0.10 | % | Moisture |
| **240-57597-13** | **F-VB-UST-13-14** | | | | | |
| 2-Butanone (MEK) | | 3.1 | J B | 21 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 60 | J | 81 | ug/Kg | 8270D |
| Di-n-butyl phthalate | | 18 | J | 81 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 13 | J B | 20 | mg/Kg | 8015C |
| Percent Solids | | 85 | | 0.10 | % | Moisture |
| Percent Moisture | | 15 | | 0.10 | % | Moisture |
| **240-57597-14** | **F-VB-UST-14-14** | | | | | |
| 2-Butanone (MEK) | | 4.0 | J B | 21 | ug/Kg | 8260C |
| Toluene | | 0.29 | J | 5.3 | ug/Kg | 8260C |
| Acetophenone | | 11 | J | 120 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 11 | J B | 19 | mg/Kg | 8015C |
| Percent Solids | | 86 | | 0.10 | % | Moisture |
| Percent Moisture | | 14 | | 0.10 | % | Moisture |
| **240-57597-15** | **F-VB-UST-DUP** | | | | | |
| 2-Butanone (MEK) | | 3.6 | J B | 21 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 22 | J | 82 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 8.2 | J B | 20 | mg/Kg | 8015C |
| Percent Solids | | 86 | | 0.10 | % | Moisture |
| Percent Moisture | | 14 | | 0.10 | % | Moisture |

# METHOD SUMMARY

Client: Tetra Tech, Inc. Job Number: 240-57597-1

| Description | Lab Location | Method | Preparation Method |
|---|---|---|---|
| **Matrix: Solid** | | | |
| Volatile Organic Compounds by GC/MS | TAL CAN | SW846 8260C | |
| Purge and Trap | TAL CAN | | SW846 5030C |
| Semivolatile Organic Compounds (GC/MS) | TAL CAN | SW846 8270D | |
| Soxhlet Extraction | TAL CAN | | SW846 3540C |
| Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics) | TAL CAN | SW846 8015C | |
| Purge and Trap | TAL CAN | | SW846 5030A |
| Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics) | TAL CAN | SW846 8015C | |
| Soxhlet Extraction | TAL CAN | | SW846 3540C |
| Polychlorinated Biphenyls (PCBs) by Gas Chromatography | TAL CAN | SW846 8082A | |
| Soxhlet Extraction | TAL CAN | | SW846 3540C |
| Percent Moisture | TAL CAN | EPA Moisture | |
| **Matrix: Water** | | | |
| Volatile Organic Compounds by GC/MS | TAL CAN | SW846 8260C | |
| Purge and Trap | TAL CAN | | SW846 5030C |

### Lab References:

TAL CAN = TestAmerica Canton

### Method References:

EPA = US Environmental Protection Agency

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

# METHOD / ANALYST SUMMARY

Client: Tetra Tech, Inc. Job Number: 240-57597-1

| Method | Analyst | Analyst ID |
|---|---|---|
| SW846 8260C | Evans, Laura | LEE |
| SW846 8260C | Lata, Todd | TJL2 |
| SW846 8260C | Macenczak, Steven | SAM |
| SW846 8270D | Gruber, John | JMG |
| SW846 8270D | Ulman, Mark | MRU |
| SW846 8015C | Reed, Rachel T | RTR |
| SW846 8015C | Bolgrin, Deborah | DEB |
| SW846 8082A | Hass, Lori | LSH |
| EPA Moisture | Nolle, Laura C | LCN |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-1-14**

Lab Sample ID: 240-57597-1 Date Sampled: 11/05/2015 0900
Client Matrix: Solid % Moisture: 15.7 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146336.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.89 g |
| Analysis Date: | 11/10/2015 1443 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.50 | U | 0.50 | 5.0 |
| 1,1,1-Trichloroethane | | 0.65 | U | 0.65 | 5.0 |
| 1,1,2,2-Tetrachloroethane | | 0.31 | U | 0.31 | 5.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.42 | U | 0.42 | 5.0 |
| 1,1-Dichloroethane | | 0.34 | U | 0.34 | 5.0 |
| 1,1-Dichloroethene | | 0.81 | U | 0.81 | 5.0 |
| 1,1-Dichloropropene | | 0.37 | U | 0.37 | 5.0 |
| 1,2,3-Trichlorobenzene | | 0.50 | U | 0.50 | 5.0 |
| 1,2,3-Trichloropropane | | 0.79 | U | 0.79 | 5.0 |
| 1,2,3-Trimethylbenzene | | 0.46 | U | 0.46 | 5.0 |
| 1,2,4-Trichlorobenzene | | 0.38 | U | 0.38 | 5.0 |
| 1,2,4-Trimethylbenzene | | 0.58 | U | 0.58 | 5.0 |
| 1,2-Dibromo-3-Chloropropane | | 2.7 | U | 2.7 | 10 |
| 1,2-Dichlorobenzene | | 0.34 | U | 0.34 | 5.0 |
| 1,2-Dichloroethane | | 0.46 | U | 0.46 | 5.0 |
| 1,2-Dichloropropane | | 0.14 | U | 0.14 | 5.0 |
| 1,3-Dichlorobenzene | | 0.69 | U | 0.69 | 5.0 |
| 1,3-Dichloropropane | | 0.40 | U | 0.40 | 5.0 |
| 1,4-Dichlorobenzene | | 0.28 | U | 0.28 | 5.0 |
| 2,2-Dichloropropane | | 0.44 | U | 0.44 | 5.0 |
| 2-Butanone (MEK) | | 3.9 | J B | 1.1 | 20 |
| 2-Chloroethyl vinyl ether | | 0.64 | U | 0.64 | 50 |
| 2-Chlorotoluene | | 0.30 | U | 0.30 | 5.0 |
| 2-Hexanone | | 0.68 | U | 0.68 | 20 |
| 4-Chlorotoluene | | 0.26 | U | 0.26 | 5.0 |
| p-Isopropyltoluene | | 0.41 | U | 0.41 | 5.0 |
| 4-Methyl-2-pentanone (MIBK) | | 1.2 | U | 1.2 | 20 |
| Acetone | | 5.7 | U | 5.7 | 20 |
| Benzene | | 0.91 | U | 0.91 | 5.0 |
| Bromobenzene | | 0.33 | U | 0.33 | 5.0 |
| Bromoform | | 0.24 | U | 0.24 | 5.0 |
| Bromomethane | | 0.42 | U | 0.42 | 5.0 |
| Carbon disulfide | | 0.59 | U | 0.59 | 5.0 |
| Carbon tetrachloride | | 0.65 | U | 0.65 | 5.0 |
| Chlorobenzene | | 0.53 | U | 0.53 | 5.0 |
| Dibromochloromethane | | 0.37 | U | 0.37 | 5.0 |
| Chloroethane | | 0.44 | U | 0.44 | 5.0 |
| Chloroform | | 0.36 | U | 0.36 | 5.0 |
| Chloromethane | | 0.80 | U | 0.80 | 5.0 |
| cis-1,2-Dichloroethene | | 0.28 | U | 0.28 | 5.0 |
| cis-1,3-Dichloropropene | | 0.72 | U | 0.72 | 5.0 |
| Dibromomethane | | 0.42 | U | 0.42 | 5.0 |
| Bromochloromethane | | 0.33 | U | 0.33 | 5.0 |
| Dichlorodifluoromethane | | 0.28 | U | 0.28 | 5.0 |
| Bromodichloromethane | | 0.29 | U | 0.29 | 5.0 |
| tert-Butyl alcohol (TBA) | | 5.8 | U | 5.8 | 200 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-1-14**

Lab Sample ID: 240-57597-1 Date Sampled: 11/05/2015 0900
Client Matrix: Solid % Moisture: 15.7 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146336.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.89 g |
| Analysis Date: | 11/10/2015 1443 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.32 | U | 0.32 | 5.0 |
| Hexachlorobutadiene | | 0.40 | U | 0.40 | 5.0 |
| Diisopropyl ether | | 0.59 | U | 0.59 | 10 |
| Isopropylbenzene | | 0.20 | U | 0.20 | 5.0 |
| Methyl tert-butyl ether | | 0.43 | U | 0.43 | 5.0 |
| Methylene Chloride | | 0.75 | U | 0.75 | 5.0 |
| m-Xylene & p-Xylene | | 0.42 | U | 0.42 | 10 |
| Naphthalene | | 0.35 | U | 0.35 | 5.0 |
| n-Butylbenzene | | 0.59 | U | 0.59 | 5.0 |
| N-Propylbenzene | | 0.60 | U | 0.60 | 5.0 |
| o-Xylene | | 0.57 | U | 0.57 | 5.0 |
| sec-Butylbenzene | | 0.92 | U | 0.92 | 5.0 |
| Styrene | | 0.40 | U | 0.40 | 5.0 |
| Tert-amyl methyl ether | | 0.43 | U | 0.43 | 5.0 |
| Ethyl t-butyl ether | | 0.41 | U | 0.41 | 5.0 |
| Tetrachloroethene | | 0.81 | U | 0.81 | 5.0 |
| Toluene | | 0.27 | U | 0.27 | 5.0 |
| trans-1,2-Dichloroethene | | 0.32 | U | 0.32 | 5.0 |
| trans-1,3-Dichloropropene | | 0.29 | U | 0.29 | 5.0 |
| Trichloroethene | | 0.38 | U | 0.38 | 5.0 |
| Trichlorofluoromethane | | 0.29 | U | 0.29 | 5.0 |
| Vinyl acetate | | 0.66 | U | 0.66 | 10 |
| Vinyl chloride | | 0.30 | U | 0.30 | 5.0 |
| Xylenes, Total | | 0.54 | U | 0.54 | 10 |
| tert-Butylbenzene | | 0.67 | U | 0.67 | 5.0 |
| 1,2-Dibromoethane | | 0.53 | U | 0.53 | 5.0 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 103 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 101 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 91 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 103 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-2-14**

Lab Sample ID: 240-57597-2 Date Sampled: 11/05/2015 0902
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146337.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.36 g |
| Analysis Date: | 11/10/2015 1505 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.55 | U | 0.55 | 5.5 |
| 1,1,1-Trichloroethane | | 0.71 | U | 0.71 | 5.5 |
| 1,1,2,2-Tetrachloroethane | | 0.34 | U | 0.34 | 5.5 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.46 | U | 0.46 | 5.5 |
| 1,1-Dichloroethane | | 0.37 | U | 0.37 | 5.5 |
| 1,1-Dichloroethene | | 0.88 | U | 0.88 | 5.5 |
| 1,1-Dichloropropene | | 0.41 | U | 0.41 | 5.5 |
| 1,2,3-Trichlorobenzene | | 0.55 | U | 0.55 | 5.5 |
| 1,2,3-Trichloropropane | | 0.86 | U | 0.86 | 5.5 |
| 1,2,3-Trimethylbenzene | | 0.50 | U | 0.50 | 5.5 |
| 1,2,4-Trichlorobenzene | | 0.42 | U | 0.42 | 5.5 |
| 1,2,4-Trimethylbenzene | | 0.64 | U | 0.64 | 5.5 |
| 1,2-Dibromo-3-Chloropropane | | 3.0 | U | 3.0 | 11 |
| 1,2-Dichlorobenzene | | 0.37 | U | 0.37 | 5.5 |
| 1,2-Dichloroethane | | 0.50 | U | 0.50 | 5.5 |
| 1,2-Dichloropropane | | 0.15 | U | 0.15 | 5.5 |
| 1,3-Dichlorobenzene | | 0.76 | U | 0.76 | 5.5 |
| 1,3-Dichloropropane | | 0.44 | U | 0.44 | 5.5 |
| 1,4-Dichlorobenzene | | 0.31 | U | 0.31 | 5.5 |
| 2,2-Dichloropropane | | 0.48 | U | 0.48 | 5.5 |
| 2-Butanone (MEK) | | 4.0 | J B | 1.2 | 22 |
| 2-Chloroethyl vinyl ether | | 0.70 | U | 0.70 | 55 |
| 2-Chlorotoluene | | 0.33 | U | 0.33 | 5.5 |
| 2-Hexanone | | 0.75 | U | 0.75 | 22 |
| 4-Chlorotoluene | | 0.29 | U | 0.29 | 5.5 |
| p-Isopropyltoluene | | 0.45 | U | 0.45 | 5.5 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 22 |
| Acetone | | 6.3 | U | 6.3 | 22 |
| Benzene | | 0.99 | U | 0.99 | 5.5 |
| Bromobenzene | | 0.36 | U | 0.36 | 5.5 |
| Bromoform | | 0.26 | U | 0.26 | 5.5 |
| Bromomethane | | 0.46 | U | 0.46 | 5.5 |
| Carbon disulfide | | 0.65 | U | 0.65 | 5.5 |
| Carbon tetrachloride | | 0.71 | U | 0.71 | 5.5 |
| Chlorobenzene | | 0.58 | U | 0.58 | 5.5 |
| Dibromochloromethane | | 0.41 | U | 0.41 | 5.5 |
| Chloroethane | | 0.48 | U | 0.48 | 5.5 |
| Chloroform | | 0.39 | U | 0.39 | 5.5 |
| Chloromethane | | 0.87 | U | 0.87 | 5.5 |
| cis-1,2-Dichloroethene | | 0.31 | U | 0.31 | 5.5 |
| cis-1,3-Dichloropropene | | 0.78 | U | 0.78 | 5.5 |
| Dibromomethane | | 0.46 | U | 0.46 | 5.5 |
| Bromochloromethane | | 0.36 | U | 0.36 | 5.5 |
| Dichlorodifluoromethane | | 0.31 | U | 0.31 | 5.5 |
| Bromodichloromethane | | 0.32 | U | 0.32 | 5.5 |
| tert-Butyl alcohol (TBA) | | 6.3 | U | 6.3 | 220 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-2-14**

Lab Sample ID: 240-57597-2 Date Sampled: 11/05/2015 0902
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146337.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.36 g |
| Analysis Date: | 11/10/2015 1505 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.35 | U | 0.35 | 5.5 |
| Hexachlorobutadiene | | 0.44 | U | 0.44 | 5.5 |
| Diisopropyl ether | | 0.65 | U | 0.65 | 11 |
| Isopropylbenzene | | 0.22 | U | 0.22 | 5.5 |
| Methyl tert-butyl ether | | 0.47 | U | 0.47 | 5.5 |
| Methylene Chloride | | 0.81 | U | 0.81 | 5.5 |
| m-Xylene & p-Xylene | | 0.46 | U | 0.46 | 11 |
| Naphthalene | | 0.38 | U | 0.38 | 5.5 |
| n-Butylbenzene | | 0.65 | U | 0.65 | 5.5 |
| N-Propylbenzene | | 0.66 | U | 0.66 | 5.5 |
| o-Xylene | | 0.63 | U | 0.63 | 5.5 |
| sec-Butylbenzene | | 1.0 | U | 1.0 | 5.5 |
| Styrene | | 0.44 | U | 0.44 | 5.5 |
| Tert-amyl methyl ether | | 0.47 | U | 0.47 | 5.5 |
| Ethyl t-butyl ether | | 0.45 | U | 0.45 | 5.5 |
| Tetrachloroethene | | 0.88 | U | 0.88 | 5.5 |
| Toluene | | 0.30 | U | 0.30 | 5.5 |
| trans-1,2-Dichloroethene | | 0.35 | U | 0.35 | 5.5 |
| trans-1,3-Dichloropropene | | 0.32 | U | 0.32 | 5.5 |
| Trichloroethene | | 0.42 | U | 0.42 | 5.5 |
| Trichlorofluoromethane | | 0.32 | U | 0.32 | 5.5 |
| Vinyl acetate | | 0.72 | U | 0.72 | 11 |
| Vinyl chloride | | 0.33 | U | 0.33 | 5.5 |
| Xylenes, Total | | 0.59 | U | 0.59 | 11 |
| tert-Butylbenzene | | 0.74 | U | 0.74 | 5.5 |
| 1,2-Dibromoethane | | 0.58 | U | 0.58 | 5.5 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 96 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 95 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 85 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 99 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-3-14**

Lab Sample ID: 240-57597-3 Date Sampled: 11/05/2015 0904
Client Matrix: Solid % Moisture: 17.8 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146338.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.23 g |
| Analysis Date: | 11/10/2015 1526 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.76 | U | 0.76 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.49 | U | 0.49 | 5.8 |
| 1,1-Dichloroethane | | 0.40 | U | 0.40 | 5.8 |
| 1,1-Dichloroethene | | 0.93 | U | 0.93 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.91 | U | 0.91 | 5.8 |
| 1,2,3-Trimethylbenzene | | 0.53 | U | 0.53 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 0.67 | U | 0.67 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.2 | U | 3.2 | 12 |
| 1,2-Dichlorobenzene | | 0.40 | U | 0.40 | 5.8 |
| 1,2-Dichloroethane | | 0.53 | U | 0.53 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.80 | U | 0.80 | 5.8 |
| 1,3-Dichloropropane | | 0.46 | U | 0.46 | 5.8 |
| 1,4-Dichlorobenzene | | 0.33 | U | 0.33 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 3.9 | J B | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.74 | U | 0.74 | 58 |
| 2-Chlorotoluene | | 0.35 | U | 0.35 | 5.8 |
| 2-Hexanone | | 0.79 | U | 0.79 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.8 |
| p-Isopropyltoluene | | 0.48 | U | 0.48 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 6.6 | U | 6.6 | 23 |
| Benzene | | 1.0 | U | 1.0 | 5.8 |
| Bromobenzene | | 0.38 | U | 0.38 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.49 | U | 0.49 | 5.8 |
| Carbon disulfide | | 0.69 | U | 0.69 | 5.8 |
| Carbon tetrachloride | | 0.76 | U | 0.76 | 5.8 |
| Chlorobenzene | | 0.62 | U | 0.62 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.42 | U | 0.42 | 5.8 |
| Chloromethane | | 0.92 | U | 0.92 | 5.8 |
| cis-1,2-Dichloroethene | | 0.33 | U | 0.33 | 5.8 |
| cis-1,3-Dichloropropene | | 0.83 | U | 0.83 | 5.8 |
| Dibromomethane | | 0.49 | U | 0.49 | 5.8 |
| Bromochloromethane | | 0.38 | U | 0.38 | 5.8 |
| Dichlorodifluoromethane | | 0.33 | U | 0.33 | 5.8 |
| Bromodichloromethane | | 0.34 | U | 0.34 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.6 | U | 6.6 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-3-14**

Lab Sample ID: 240-57597-3 Date Sampled: 11/05/2015 0904
Client Matrix: Solid % Moisture: 17.8 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146338.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.23 g |
| Analysis Date: | 11/10/2015 1526 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.37 | U | 0.37 | 5.8 |
| Hexachlorobutadiene | | 0.46 | U | 0.46 | 5.8 |
| Diisopropyl ether | | 0.69 | U | 0.69 | 12 |
| Isopropylbenzene | | 0.23 | U | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.86 | U | 0.86 | 5.8 |
| m-Xylene & p-Xylene | | 0.49 | U | 0.49 | 12 |
| Naphthalene | | 0.41 | U | 0.41 | 5.8 |
| n-Butylbenzene | | 0.69 | U | 0.69 | 5.8 |
| N-Propylbenzene | | 0.70 | U | 0.70 | 5.8 |
| o-Xylene | | 0.66 | U | 0.66 | 5.8 |
| sec-Butylbenzene | | 1.1 | U | 1.1 | 5.8 |
| Styrene | | 0.46 | U | 0.46 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.48 | U | 0.48 | 5.8 |
| Tetrachloroethene | | 0.93 | U | 0.93 | 5.8 |
| Toluene | | 0.31 | U | 0.31 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.34 | U | 0.34 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.34 | U | 0.34 | 5.8 |
| Vinyl acetate | | 0.77 | U | 0.77 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 0.63 | U | 0.63 | 12 |
| tert-Butylbenzene | | 0.78 | U | 0.78 | 5.8 |
| 1,2-Dibromoethane | | 0.62 | U | 0.62 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 93 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 90 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 86 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 90 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-4-14**

Lab Sample ID: 240-57597-4 Date Sampled: 11/05/2015 0906
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146339.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.43 g |
| Analysis Date: | 11/10/2015 1548 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.55 | U | 0.55 | 5.5 |
| 1,1,1-Trichloroethane | | 0.71 | U | 0.71 | 5.5 |
| 1,1,2,2-Tetrachloroethane | | 0.34 | U | 0.34 | 5.5 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.46 | U | 0.46 | 5.5 |
| 1,1-Dichloroethane | | 0.37 | U | 0.37 | 5.5 |
| 1,1-Dichloroethene | | 0.87 | U | 0.87 | 5.5 |
| 1,1-Dichloropropene | | 0.40 | U | 0.40 | 5.5 |
| 1,2,3-Trichlorobenzene | | 0.55 | U | 0.55 | 5.5 |
| 1,2,3-Trichloropropane | | 0.85 | U | 0.85 | 5.5 |
| 1,2,3-Trimethylbenzene | | 0.50 | U | 0.50 | 5.5 |
| 1,2,4-Trichlorobenzene | | 0.41 | U | 0.41 | 5.5 |
| 1,2,4-Trimethylbenzene | | 0.63 | U | 0.63 | 5.5 |
| 1,2-Dibromo-3-Chloropropane | | 3.0 | U | 3.0 | 11 |
| 1,2-Dichlorobenzene | | 0.37 | U | 0.37 | 5.5 |
| 1,2-Dichloroethane | | 0.50 | U | 0.50 | 5.5 |
| 1,2-Dichloropropane | | 0.15 | U | 0.15 | 5.5 |
| 1,3-Dichlorobenzene | | 0.75 | U | 0.75 | 5.5 |
| 1,3-Dichloropropane | | 0.44 | U | 0.44 | 5.5 |
| 1,4-Dichlorobenzene | | 0.31 | U | 0.31 | 5.5 |
| 2,2-Dichloropropane | | 0.48 | U | 0.48 | 5.5 |
| 2-Butanone (MEK) | | 4.0 | J B | 1.2 | 22 |
| 2-Chloroethyl vinyl ether | | 0.70 | U | 0.70 | 55 |
| 2-Chlorotoluene | | 0.33 | U | 0.33 | 5.5 |
| 2-Hexanone | | 0.74 | U | 0.74 | 22 |
| 4-Chlorotoluene | | 0.28 | U | 0.28 | 5.5 |
| p-Isopropyltoluene | | 0.45 | U | 0.45 | 5.5 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 22 |
| Acetone | | 6.2 | U | 6.2 | 22 |
| Benzene | | 0.98 | U | 0.98 | 5.5 |
| Bromobenzene | | 0.36 | U | 0.36 | 5.5 |
| Bromoform | | 0.26 | U | 0.26 | 5.5 |
| Bromomethane | | 0.46 | U | 0.46 | 5.5 |
| Carbon disulfide | | 0.64 | U | 0.64 | 5.5 |
| Carbon tetrachloride | | 0.71 | U | 0.71 | 5.5 |
| Chlorobenzene | | 0.58 | U | 0.58 | 5.5 |
| Dibromochloromethane | | 0.40 | U | 0.40 | 5.5 |
| Chloroethane | | 0.48 | U | 0.48 | 5.5 |
| Chloroform | | 0.39 | U | 0.39 | 5.5 |
| Chloromethane | | 0.86 | U | 0.86 | 5.5 |
| cis-1,2-Dichloroethene | | 0.31 | U | 0.31 | 5.5 |
| cis-1,3-Dichloropropene | | 0.77 | U | 0.77 | 5.5 |
| Dibromomethane | | 0.46 | U | 0.46 | 5.5 |
| Bromochloromethane | | 0.36 | U | 0.36 | 5.5 |
| Dichlorodifluoromethane | | 0.31 | U | 0.31 | 5.5 |
| Bromodichloromethane | | 0.32 | U | 0.32 | 5.5 |
| tert-Butyl alcohol (TBA) | | 6.2 | U | 6.2 | 220 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-4-14**

Lab Sample ID: 240-57597-4 Date Sampled: 11/05/2015 0906
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146339.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.43 g |
| Analysis Date: | 11/10/2015 1548 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.35 | U | 0.35 | 5.5 |
| Hexachlorobutadiene | | 0.44 | U | 0.44 | 5.5 |
| Diisopropyl ether | | 0.64 | U | 0.64 | 11 |
| Isopropylbenzene | | 0.22 | U | 0.22 | 5.5 |
| Methyl tert-butyl ether | | 0.47 | U | 0.47 | 5.5 |
| Methylene Chloride | | 0.81 | U | 0.81 | 5.5 |
| m-Xylene & p-Xylene | | 0.46 | U | 0.46 | 11 |
| Naphthalene | | 0.38 | U | 0.38 | 5.5 |
| n-Butylbenzene | | 0.64 | U | 0.64 | 5.5 |
| N-Propylbenzene | | 0.65 | U | 0.65 | 5.5 |
| o-Xylene | | 0.62 | U | 0.62 | 5.5 |
| sec-Butylbenzene | | 0.99 | U | 0.99 | 5.5 |
| Styrene | | 0.44 | U | 0.44 | 5.5 |
| Tert-amyl methyl ether | | 0.47 | U | 0.47 | 5.5 |
| Ethyl t-butyl ether | | 0.45 | U | 0.45 | 5.5 |
| Tetrachloroethene | | 0.87 | U | 0.87 | 5.5 |
| Toluene | | 0.29 | U | 0.29 | 5.5 |
| trans-1,2-Dichloroethene | | 0.35 | U | 0.35 | 5.5 |
| trans-1,3-Dichloropropene | | 0.32 | U | 0.32 | 5.5 |
| Trichloroethene | | 0.41 | U | 0.41 | 5.5 |
| Trichlorofluoromethane | | 0.32 | U | 0.32 | 5.5 |
| Vinyl acetate | | 0.72 | U | 0.72 | 11 |
| Vinyl chloride | | 0.33 | U | 0.33 | 5.5 |
| Xylenes, Total | | 0.59 | U | 0.59 | 11 |
| tert-Butylbenzene | | 0.73 | U | 0.73 | 5.5 |
| 1,2-Dibromoethane | | 0.58 | U | 0.58 | 5.5 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 93 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 94 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 83 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 102 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-5-14**

Lab Sample ID: 240-57597-5 Date Sampled: 11/05/2015 0908
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146340.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.09 g |
| Analysis Date: | 11/10/2015 1609 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.75 | U | 0.75 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.49 | U | 0.49 | 5.8 |
| 1,1-Dichloroethane | | 0.39 | U | 0.39 | 5.8 |
| 1,1-Dichloroethene | | 0.93 | U | 0.93 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.90 | U | 0.90 | 5.8 |
| 1,2,3-Trimethylbenzene | | 0.53 | U | 0.53 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U F2 | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 0.67 | U | 0.67 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.1 | U | 3.1 | 12 |
| 1,2-Dichlorobenzene | | 0.39 | U | 0.39 | 5.8 |
| 1,2-Dichloroethane | | 0.53 | U | 0.53 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.80 | U | 0.80 | 5.8 |
| 1,3-Dichloropropane | | 0.46 | U | 0.46 | 5.8 |
| 1,4-Dichlorobenzene | | 0.32 | U | 0.32 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 4.3 | J B | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.74 | U | 0.74 | 58 |
| 2-Chlorotoluene | | 0.35 | U | 0.35 | 5.8 |
| 2-Hexanone | | 0.79 | U | 0.79 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.8 |
| p-Isopropyltoluene | | 0.47 | U F2 | 0.47 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U F1 | 1.4 | 23 |
| Acetone | | 13 | J | 6.6 | 23 |
| Benzene | | 1.0 | U | 1.0 | 5.8 |
| Bromobenzene | | 0.38 | U | 0.38 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.49 | U | 0.49 | 5.8 |
| Carbon disulfide | | 0.68 | U | 0.68 | 5.8 |
| Carbon tetrachloride | | 0.75 | U | 0.75 | 5.8 |
| Chlorobenzene | | 0.61 | U | 0.61 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.42 | U | 0.42 | 5.8 |
| Chloromethane | | 0.92 | U | 0.92 | 5.8 |
| cis-1,2-Dichloroethene | | 0.32 | U | 0.32 | 5.8 |
| cis-1,3-Dichloropropene | | 0.82 | U | 0.82 | 5.8 |
| Dibromomethane | | 0.49 | U | 0.49 | 5.8 |
| Bromochloromethane | | 0.38 | U | 0.38 | 5.8 |
| Dichlorodifluoromethane | | 0.32 | U | 0.32 | 5.8 |
| Bromodichloromethane | | 0.34 | U | 0.34 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.6 | U | 6.6 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-5-14**

Lab Sample ID: 240-57597-5 Date Sampled: 11/05/2015 0908
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146340.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.09 g |
| Analysis Date: | 11/10/2015 1609 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.37 | U | 0.37 | 5.8 |
| Hexachlorobutadiene | | 0.46 | U F2 | 0.46 | 5.8 |
| Diisopropyl ether | | 0.68 | U | 0.68 | 12 |
| Isopropylbenzene | | 0.23 | U | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.86 | U | 0.86 | 5.8 |
| m-Xylene & p-Xylene | | 0.49 | U | 0.49 | 12 |
| Naphthalene | | 0.41 | J | 0.41 | 5.8 |
| n-Butylbenzene | | 0.68 | U F2 | 0.68 | 5.8 |
| N-Propylbenzene | | 0.70 | U | 0.70 | 5.8 |
| o-Xylene | | 0.66 | U | 0.66 | 5.8 |
| sec-Butylbenzene | | 1.1 | U F2 | 1.1 | 5.8 |
| Styrene | | 0.46 | U | 0.46 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.47 | U | 0.47 | 5.8 |
| Tetrachloroethene | | 0.93 | U | 0.93 | 5.8 |
| Toluene | | 0.31 | U | 0.31 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.34 | U | 0.34 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.34 | U | 0.34 | 5.8 |
| Vinyl acetate | | 0.76 | U F1 | 0.76 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 0.63 | U | 0.63 | 12 |
| tert-Butylbenzene | | 0.78 | U | 0.78 | 5.8 |
| 1,2-Dibromoethane | | 0.61 | U | 0.61 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 100 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 101 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 87 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 104 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-6-14**

Lab Sample ID: 240-57597-6 Date Sampled: 11/05/2015 0910
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146343.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.18 g |
| Analysis Date: | 11/10/2015 1713 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.75 | U | 0.75 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.49 | U | 0.49 | 5.8 |
| 1,1-Dichloroethane | | 0.39 | U | 0.39 | 5.8 |
| 1,1-Dichloroethene | | 0.93 | U | 0.93 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.90 | U | 0.90 | 5.8 |
| 1,2,3-Trimethylbenzene | | 0.53 | U | 0.53 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 0.67 | U | 0.67 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.1 | U | 3.1 | 12 |
| 1,2-Dichlorobenzene | | 0.39 | U | 0.39 | 5.8 |
| 1,2-Dichloroethane | | 0.53 | U | 0.53 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.80 | U | 0.80 | 5.8 |
| 1,3-Dichloropropane | | 0.46 | U | 0.46 | 5.8 |
| 1,4-Dichlorobenzene | | 0.32 | U | 0.32 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 3.8 | J B | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.74 | U | 0.74 | 58 |
| 2-Chlorotoluene | | 0.35 | U | 0.35 | 5.8 |
| 2-Hexanone | | 0.79 | U | 0.79 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.8 |
| p-Isopropyltoluene | | 0.47 | U | 0.47 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 6.6 | U | 6.6 | 23 |
| Benzene | | 1.0 | U | 1.0 | 5.8 |
| Bromobenzene | | 0.38 | U | 0.38 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.49 | U | 0.49 | 5.8 |
| Carbon disulfide | | 0.68 | U | 0.68 | 5.8 |
| Carbon tetrachloride | | 0.75 | U | 0.75 | 5.8 |
| Chlorobenzene | | 0.61 | U | 0.61 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.42 | U | 0.42 | 5.8 |
| Chloromethane | | 0.91 | U | 0.91 | 5.8 |
| cis-1,2-Dichloroethene | | 0.32 | U | 0.32 | 5.8 |
| cis-1,3-Dichloropropene | | 0.82 | U | 0.82 | 5.8 |
| Dibromomethane | | 0.49 | U | 0.49 | 5.8 |
| Bromochloromethane | | 0.38 | U | 0.38 | 5.8 |
| Dichlorodifluoromethane | | 0.32 | U | 0.32 | 5.8 |
| Bromodichloromethane | | 0.34 | U | 0.34 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.6 | U | 6.6 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-6-14**

Lab Sample ID: 240-57597-6 Date Sampled: 11/05/2015 0910
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146343.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.18 g |
| Analysis Date: | 11/10/2015 1713 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.37 | U | 0.37 | 5.8 |
| Hexachlorobutadiene | | 0.46 | U | 0.46 | 5.8 |
| Diisopropyl ether | | 0.68 | U | 0.68 | 12 |
| Isopropylbenzene | | 0.23 | U | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.86 | U | 0.86 | 5.8 |
| m-Xylene & p-Xylene | | 0.49 | U | 0.49 | 12 |
| Naphthalene | | 0.45 | J | 0.40 | 5.8 |
| n-Butylbenzene | | 0.68 | U | 0.68 | 5.8 |
| N-Propylbenzene | | 0.69 | U | 0.69 | 5.8 |
| o-Xylene | | 0.66 | U | 0.66 | 5.8 |
| sec-Butylbenzene | | 1.1 | U | 1.1 | 5.8 |
| Styrene | | 0.46 | U | 0.46 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.47 | U | 0.47 | 5.8 |
| Tetrachloroethene | | 0.93 | U | 0.93 | 5.8 |
| Toluene | | 0.31 | U | 0.31 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.34 | U | 0.34 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.34 | U | 0.34 | 5.8 |
| Vinyl acetate | | 0.76 | U | 0.76 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 0.62 | U | 0.62 | 12 |
| tert-Butylbenzene | | 0.77 | U | 0.77 | 5.8 |
| 1,2-Dibromoethane | | 0.61 | U | 0.61 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 88 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 89 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 81 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 93 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-7-14**

Lab Sample ID: 240-57597-7 Date Sampled: 11/05/2015 0912
Client Matrix: Solid % Moisture: 16.2 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146344.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.31 g |
| Analysis Date: | 11/10/2015 1735 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.56 | U | 0.56 | 5.6 |
| 1,1,1-Trichloroethane | | 0.73 | U | 0.73 | 5.6 |
| 1,1,2,2-Tetrachloroethane | | 0.35 | U | 0.35 | 5.6 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.47 | U | 0.47 | 5.6 |
| 1,1-Dichloroethane | | 0.38 | U | 0.38 | 5.6 |
| 1,1-Dichloroethene | | 0.90 | U | 0.90 | 5.6 |
| 1,1-Dichloropropene | | 0.42 | U | 0.42 | 5.6 |
| 1,2,3-Trichlorobenzene | | 0.56 | U | 0.56 | 5.6 |
| 1,2,3-Trichloropropane | | 0.88 | U | 0.88 | 5.6 |
| 1,2,3-Trimethylbenzene | | 0.52 | U | 0.52 | 5.6 |
| 1,2,4-Trichlorobenzene | | 0.43 | U | 0.43 | 5.6 |
| 1,2,4-Trimethylbenzene | | 0.65 | U | 0.65 | 5.6 |
| 1,2-Dibromo-3-Chloropropane | | 3.0 | U | 3.0 | 11 |
| 1,2-Dichlorobenzene | | 0.38 | U | 0.38 | 5.6 |
| 1,2-Dichloroethane | | 0.52 | U | 0.52 | 5.6 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.6 |
| 1,3-Dichlorobenzene | | 0.78 | U | 0.78 | 5.6 |
| 1,3-Dichloropropane | | 0.45 | U | 0.45 | 5.6 |
| 1,4-Dichlorobenzene | | 0.31 | U | 0.31 | 5.6 |
| 2,2-Dichloropropane | | 0.49 | U | 0.49 | 5.6 |
| 2-Butanone (MEK) | | 4.1 | J B | 1.3 | 22 |
| 2-Chloroethyl vinyl ether | | 0.72 | U | 0.72 | 56 |
| 2-Chlorotoluene | | 0.34 | U | 0.34 | 5.6 |
| 2-Hexanone | | 0.76 | U | 0.76 | 22 |
| 4-Chlorotoluene | | 0.29 | U | 0.29 | 5.6 |
| p-Isopropyltoluene | | 0.46 | U | 0.46 | 5.6 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 22 |
| Acetone | | 6.4 | U | 6.4 | 22 |
| Benzene | | 1.0 | U | 1.0 | 5.6 |
| Bromobenzene | | 0.37 | U | 0.37 | 5.6 |
| Bromoform | | 0.27 | U | 0.27 | 5.6 |
| Bromomethane | | 0.47 | U | 0.47 | 5.6 |
| Carbon disulfide | | 0.66 | U | 0.66 | 5.6 |
| Carbon tetrachloride | | 0.73 | U | 0.73 | 5.6 |
| Chlorobenzene | | 0.60 | U | 0.60 | 5.6 |
| Dibromochloromethane | | 0.42 | U | 0.42 | 5.6 |
| Chloroethane | | 0.49 | U | 0.49 | 5.6 |
| Chloroform | | 0.40 | U | 0.40 | 5.6 |
| Chloromethane | | 0.89 | U | 0.89 | 5.6 |
| cis-1,2-Dichloroethene | | 0.31 | U | 0.31 | 5.6 |
| cis-1,3-Dichloropropene | | 0.80 | U | 0.80 | 5.6 |
| Dibromomethane | | 0.47 | U | 0.47 | 5.6 |
| Bromochloromethane | | 0.37 | U | 0.37 | 5.6 |
| Dichlorodifluoromethane | | 0.31 | U | 0.31 | 5.6 |
| Bromodichloromethane | | 0.33 | U | 0.33 | 5.6 |
| tert-Butyl alcohol (TBA) | | 6.4 | U | 6.4 | 220 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-7-14**

Lab Sample ID: 240-57597-7 Date Sampled: 11/05/2015 0912
Client Matrix: Solid % Moisture: 16.2 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146344.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.31 g |
| Analysis Date: | 11/10/2015 1735 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.36 | U | 0.36 | 5.6 |
| Hexachlorobutadiene | | 0.45 | U | 0.45 | 5.6 |
| Diisopropyl ether | | 0.66 | U | 0.66 | 11 |
| Isopropylbenzene | | 0.22 | U | 0.22 | 5.6 |
| Methyl tert-butyl ether | | 0.48 | U | 0.48 | 5.6 |
| Methylene Chloride | | 0.83 | U | 0.83 | 5.6 |
| m-Xylene & p-Xylene | | 0.47 | U | 0.47 | 11 |
| Naphthalene | | 0.39 | U | 0.39 | 5.6 |
| n-Butylbenzene | | 0.66 | U | 0.66 | 5.6 |
| N-Propylbenzene | | 0.67 | U | 0.67 | 5.6 |
| o-Xylene | | 0.64 | U | 0.64 | 5.6 |
| sec-Butylbenzene | | 1.0 | U | 1.0 | 5.6 |
| Styrene | | 0.45 | U | 0.45 | 5.6 |
| Tert-amyl methyl ether | | 0.48 | U | 0.48 | 5.6 |
| Ethyl t-butyl ether | | 0.46 | U | 0.46 | 5.6 |
| Tetrachloroethene | | 0.90 | U | 0.90 | 5.6 |
| Toluene | | 0.30 | U | 0.30 | 5.6 |
| trans-1,2-Dichloroethene | | 0.36 | U | 0.36 | 5.6 |
| trans-1,3-Dichloropropene | | 0.33 | U | 0.33 | 5.6 |
| Trichloroethene | | 0.43 | U | 0.43 | 5.6 |
| Trichlorofluoromethane | | 0.33 | U | 0.33 | 5.6 |
| Vinyl acetate | | 0.74 | U | 0.74 | 11 |
| Vinyl chloride | | 0.34 | U | 0.34 | 5.6 |
| Xylenes, Total | | 0.61 | U | 0.61 | 11 |
| tert-Butylbenzene | | 0.75 | U | 0.75 | 5.6 |
| 1,2-Dibromoethane | | 0.60 | U | 0.60 | 5.6 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 98 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 98 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 87 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 103 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-8-14**

Lab Sample ID: 240-57597-8 Date Sampled: 11/05/2015 0914
Client Matrix: Solid % Moisture: 18.0 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146345.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.29 g |
| Analysis Date: | 11/10/2015 1756 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.75 | U | 0.75 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.48 | U | 0.48 | 5.8 |
| 1,1-Dichloroethane | | 0.39 | U | 0.39 | 5.8 |
| 1,1-Dichloroethene | | 0.92 | U | 0.92 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.90 | U | 0.90 | 5.8 |
| 1,2,3-Trimethylbenzene | | 0.53 | U | 0.53 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 0.67 | U | 0.67 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.1 | U | 3.1 | 12 |
| 1,2-Dichlorobenzene | | 0.39 | U | 0.39 | 5.8 |
| 1,2-Dichloroethane | | 0.53 | U | 0.53 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.80 | U | 0.80 | 5.8 |
| 1,3-Dichloropropane | | 0.46 | U | 0.46 | 5.8 |
| 1,4-Dichlorobenzene | | 0.32 | U | 0.32 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 4.3 | J B | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.74 | U | 0.74 | 58 |
| 2-Chlorotoluene | | 0.35 | U | 0.35 | 5.8 |
| 2-Hexanone | | 0.78 | U | 0.78 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.8 |
| p-Isopropyltoluene | | 0.47 | U | 0.47 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 6.6 | U | 6.6 | 23 |
| Benzene | | 1.0 | U | 1.0 | 5.8 |
| Bromobenzene | | 0.38 | U | 0.38 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.48 | U | 0.48 | 5.8 |
| Carbon disulfide | | 0.68 | U | 0.68 | 5.8 |
| Carbon tetrachloride | | 0.75 | U | 0.75 | 5.8 |
| Chlorobenzene | | 0.61 | U | 0.61 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.41 | U | 0.41 | 5.8 |
| Chloromethane | | 0.91 | U | 0.91 | 5.8 |
| cis-1,2-Dichloroethene | | 0.32 | U | 0.32 | 5.8 |
| cis-1,3-Dichloropropene | | 0.82 | U | 0.82 | 5.8 |
| Dibromomethane | | 0.48 | U | 0.48 | 5.8 |
| Bromochloromethane | | 0.38 | U | 0.38 | 5.8 |
| Dichlorodifluoromethane | | 0.32 | U | 0.32 | 5.8 |
| Bromodichloromethane | | 0.33 | U | 0.33 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.6 | U | 6.6 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-8-14**

Lab Sample ID: 240-57597-8 Date Sampled: 11/05/2015 0914
Client Matrix: Solid % Moisture: 18.0 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146345.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.29 g |
| Analysis Date: | 11/10/2015 1756 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.37 | U | 0.37 | 5.8 |
| Hexachlorobutadiene | | 0.46 | U | 0.46 | 5.8 |
| Diisopropyl ether | | 0.68 | U | 0.68 | 12 |
| Isopropylbenzene | | 0.23 | U | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.85 | U | 0.85 | 5.8 |
| m-Xylene & p-Xylene | | 0.48 | U | 0.48 | 12 |
| Naphthalene | | 0.40 | U | 0.40 | 5.8 |
| n-Butylbenzene | | 0.68 | U | 0.68 | 5.8 |
| N-Propylbenzene | | 0.69 | U | 0.69 | 5.8 |
| o-Xylene | | 0.66 | U | 0.66 | 5.8 |
| sec-Butylbenzene | | 1.0 | U | 1.0 | 5.8 |
| Styrene | | 0.46 | U | 0.46 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.47 | U | 0.47 | 5.8 |
| Tetrachloroethene | | 0.92 | U | 0.92 | 5.8 |
| Toluene | | 0.33 | J | 0.31 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.33 | U | 0.33 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.33 | U | 0.33 | 5.8 |
| Vinyl acetate | | 0.76 | U | 0.76 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 0.62 | U | 0.62 | 12 |
| tert-Butylbenzene | | 0.77 | U | 0.77 | 5.8 |
| 1,2-Dibromoethane | | 0.61 | U | 0.61 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 97 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 93 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 85 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 93 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-9-14**

Lab Sample ID: 240-57597-9 Date Sampled: 11/05/2015 0916
Client Matrix: Solid % Moisture: 19.1 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146346.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.35 g |
| Analysis Date: | 11/10/2015 1818 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.75 | U | 0.75 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.49 | U | 0.49 | 5.8 |
| 1,1-Dichloroethane | | 0.39 | U | 0.39 | 5.8 |
| 1,1-Dichloroethene | | 0.92 | U | 0.92 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.90 | U | 0.90 | 5.8 |
| 1,2,3-Trimethylbenzene | | 0.53 | U | 0.53 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 0.67 | U | 0.67 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.1 | U | 3.1 | 12 |
| 1,2-Dichlorobenzene | | 0.39 | U | 0.39 | 5.8 |
| 1,2-Dichloroethane | | 0.53 | U | 0.53 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.80 | U | 0.80 | 5.8 |
| 1,3-Dichloropropane | | 0.46 | U | 0.46 | 5.8 |
| 1,4-Dichlorobenzene | | 0.32 | U | 0.32 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 4.4 | J B | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.74 | U | 0.74 | 58 |
| 2-Chlorotoluene | | 0.35 | U | 0.35 | 5.8 |
| 2-Hexanone | | 0.79 | U | 0.79 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.8 |
| p-Isopropyltoluene | | 0.47 | U | 0.47 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 12 | J | 6.6 | 23 |
| Benzene | | 1.0 | U | 1.0 | 5.8 |
| Bromobenzene | | 0.38 | U | 0.38 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.49 | U | 0.49 | 5.8 |
| Carbon disulfide | | 0.68 | U | 0.68 | 5.8 |
| Carbon tetrachloride | | 0.75 | U | 0.75 | 5.8 |
| Chlorobenzene | | 0.61 | U | 0.61 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.42 | U | 0.42 | 5.8 |
| Chloromethane | | 0.91 | U | 0.91 | 5.8 |
| cis-1,2-Dichloroethene | | 0.32 | U | 0.32 | 5.8 |
| cis-1,3-Dichloropropene | | 0.82 | U | 0.82 | 5.8 |
| Dibromomethane | | 0.49 | U | 0.49 | 5.8 |
| Bromochloromethane | | 0.38 | U | 0.38 | 5.8 |
| Dichlorodifluoromethane | | 0.32 | U | 0.32 | 5.8 |
| Bromodichloromethane | | 0.34 | U | 0.34 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.6 | U | 6.6 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-9-14**

Lab Sample ID: 240-57597-9 Date Sampled: 11/05/2015 0916
Client Matrix: Solid % Moisture: 19.1 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146346.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.35 g |
| Analysis Date: | 11/10/2015 1818 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.37 | U | 0.37 | 5.8 |
| Hexachlorobutadiene | | 0.46 | U | 0.46 | 5.8 |
| Diisopropyl ether | | 0.68 | U | 0.68 | 12 |
| Isopropylbenzene | | 0.23 | U | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.85 | U | 0.85 | 5.8 |
| m-Xylene & p-Xylene | | 0.49 | U | 0.49 | 12 |
| Naphthalene | | 0.40 | U | 0.40 | 5.8 |
| n-Butylbenzene | | 0.68 | U | 0.68 | 5.8 |
| N-Propylbenzene | | 0.69 | U | 0.69 | 5.8 |
| o-Xylene | | 0.66 | U | 0.66 | 5.8 |
| sec-Butylbenzene | | 1.1 | U | 1.1 | 5.8 |
| Styrene | | 0.46 | U | 0.46 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.47 | U | 0.47 | 5.8 |
| Tetrachloroethene | | 0.92 | U | 0.92 | 5.8 |
| Toluene | | 0.36 | J | 0.31 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.34 | U | 0.34 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.34 | U | 0.34 | 5.8 |
| Vinyl acetate | | 0.76 | U | 0.76 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 0.62 | U | 0.62 | 12 |
| tert-Butylbenzene | | 0.77 | U | 0.77 | 5.8 |
| 1,2-Dibromoethane | | 0.61 | U | 0.61 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 89 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 88 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 82 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 94 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-10-14**

Lab Sample ID: 240-57597-10 Date Sampled: 11/05/2015 0918
Client Matrix: Solid % Moisture: 13.4 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146347.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.14 g |
| Analysis Date: | 11/10/2015 1839 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.56 | U | 0.56 | 5.6 |
| 1,1,1-Trichloroethane | | 0.73 | U | 0.73 | 5.6 |
| 1,1,2,2-Tetrachloroethane | | 0.35 | U | 0.35 | 5.6 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.47 | U | 0.47 | 5.6 |
| 1,1-Dichloroethane | | 0.38 | U | 0.38 | 5.6 |
| 1,1-Dichloroethene | | 0.90 | U | 0.90 | 5.6 |
| 1,1-Dichloropropene | | 0.42 | U | 0.42 | 5.6 |
| 1,2,3-Trichlorobenzene | | 0.56 | U | 0.56 | 5.6 |
| 1,2,3-Trichloropropane | | 0.88 | U | 0.88 | 5.6 |
| 1,2,3-Trimethylbenzene | | 0.52 | U | 0.52 | 5.6 |
| 1,2,4-Trichlorobenzene | | 0.43 | U | 0.43 | 5.6 |
| 1,2,4-Trimethylbenzene | | 0.65 | U | 0.65 | 5.6 |
| 1,2-Dibromo-3-Chloropropane | | 3.0 | U | 3.0 | 11 |
| 1,2-Dichlorobenzene | | 0.38 | U | 0.38 | 5.6 |
| 1,2-Dichloroethane | | 0.52 | U | 0.52 | 5.6 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.6 |
| 1,3-Dichlorobenzene | | 0.78 | U | 0.78 | 5.6 |
| 1,3-Dichloropropane | | 0.45 | U | 0.45 | 5.6 |
| 1,4-Dichlorobenzene | | 0.31 | U | 0.31 | 5.6 |
| 2,2-Dichloropropane | | 0.49 | U | 0.49 | 5.6 |
| 2-Butanone (MEK) | | 3.5 | J B | 1.3 | 22 |
| 2-Chloroethyl vinyl ether | | 0.72 | U | 0.72 | 56 |
| 2-Chlorotoluene | | 0.34 | U | 0.34 | 5.6 |
| 2-Hexanone | | 0.76 | U | 0.76 | 22 |
| 4-Chlorotoluene | | 0.29 | U | 0.29 | 5.6 |
| p-Isopropyltoluene | | 0.46 | U | 0.46 | 5.6 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 22 |
| Acetone | | 6.4 | U | 6.4 | 22 |
| Benzene | | 1.0 | U | 1.0 | 5.6 |
| Bromobenzene | | 0.37 | U | 0.37 | 5.6 |
| Bromoform | | 0.27 | U | 0.27 | 5.6 |
| Bromomethane | | 0.47 | U | 0.47 | 5.6 |
| Carbon disulfide | | 0.66 | U | 0.66 | 5.6 |
| Carbon tetrachloride | | 0.73 | U | 0.73 | 5.6 |
| Chlorobenzene | | 0.60 | U | 0.60 | 5.6 |
| Dibromochloromethane | | 0.42 | U | 0.42 | 5.6 |
| Chloroethane | | 0.49 | U | 0.49 | 5.6 |
| Chloroform | | 0.40 | U | 0.40 | 5.6 |
| Chloromethane | | 0.89 | U | 0.89 | 5.6 |
| cis-1,2-Dichloroethene | | 0.31 | U | 0.31 | 5.6 |
| cis-1,3-Dichloropropene | | 0.80 | U | 0.80 | 5.6 |
| Dibromomethane | | 0.47 | U | 0.47 | 5.6 |
| Bromochloromethane | | 0.37 | U | 0.37 | 5.6 |
| Dichlorodifluoromethane | | 0.31 | U | 0.31 | 5.6 |
| Bromodichloromethane | | 0.33 | U | 0.33 | 5.6 |
| tert-Butyl alcohol (TBA) | | 6.4 | U | 6.4 | 220 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-10-14**

Lab Sample ID: 240-57597-10 Date Sampled: 11/05/2015 0918
Client Matrix: Solid % Moisture: 13.4 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146347.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.14 g |
| Analysis Date: | 11/10/2015 1839 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.36 | U | 0.36 | 5.6 |
| Hexachlorobutadiene | | 0.45 | U | 0.45 | 5.6 |
| Diisopropyl ether | | 0.66 | U | 0.66 | 11 |
| Isopropylbenzene | | 0.22 | U | 0.22 | 5.6 |
| Methyl tert-butyl ether | | 0.48 | U | 0.48 | 5.6 |
| Methylene Chloride | | 0.83 | U | 0.83 | 5.6 |
| m-Xylene & p-Xylene | | 0.47 | U | 0.47 | 11 |
| Naphthalene | | 0.39 | U | 0.39 | 5.6 |
| n-Butylbenzene | | 0.66 | U | 0.66 | 5.6 |
| N-Propylbenzene | | 0.67 | U | 0.67 | 5.6 |
| o-Xylene | | 0.64 | U | 0.64 | 5.6 |
| sec-Butylbenzene | | 1.0 | U | 1.0 | 5.6 |
| Styrene | | 0.45 | U | 0.45 | 5.6 |
| Tert-amyl methyl ether | | 0.48 | U | 0.48 | 5.6 |
| Ethyl t-butyl ether | | 0.46 | U | 0.46 | 5.6 |
| Tetrachloroethene | | 0.90 | U | 0.90 | 5.6 |
| Toluene | | 0.30 | U | 0.30 | 5.6 |
| trans-1,2-Dichloroethene | | 0.36 | U | 0.36 | 5.6 |
| trans-1,3-Dichloropropene | | 0.33 | U | 0.33 | 5.6 |
| Trichloroethene | | 0.43 | U | 0.43 | 5.6 |
| Trichlorofluoromethane | | 0.33 | U | 0.33 | 5.6 |
| Vinyl acetate | | 0.74 | U | 0.74 | 11 |
| Vinyl chloride | | 0.34 | U | 0.34 | 5.6 |
| Xylenes, Total | | 0.61 | U | 0.61 | 11 |
| tert-Butylbenzene | | 0.75 | U | 0.75 | 5.6 |
| 1,2-Dibromoethane | | 0.60 | U | 0.60 | 5.6 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 89 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 88 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 83 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 93 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-11-14**

Lab Sample ID: 240-57597-11 Date Sampled: 11/05/2015 0920
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146348.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.03 g |
| Analysis Date: | 11/10/2015 1901 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.60 | U | 0.60 | 6.0 |
| 1,1,1-Trichloroethane | | 0.77 | U | 0.77 | 6.0 |
| 1,1,2,2-Tetrachloroethane | | 0.37 | U | 0.37 | 6.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.50 | U | 0.50 | 6.0 |
| 1,1-Dichloroethane | | 0.40 | U | 0.40 | 6.0 |
| 1,1-Dichloroethene | | 0.95 | U | 0.95 | 6.0 |
| 1,1-Dichloropropene | | 0.44 | U | 0.44 | 6.0 |
| 1,2,3-Trichlorobenzene | | 0.60 | U | 0.60 | 6.0 |
| 1,2,3-Trichloropropane | | 0.93 | U | 0.93 | 6.0 |
| 1,2,3-Trimethylbenzene | | 0.55 | U | 0.55 | 6.0 |
| 1,2,4-Trichlorobenzene | | 0.45 | U | 0.45 | 6.0 |
| 1,2,4-Trimethylbenzene | | 0.69 | U | 0.69 | 6.0 |
| 1,2-Dibromo-3-Chloropropane | | 3.2 | U | 3.2 | 12 |
| 1,2-Dichlorobenzene | | 0.40 | U | 0.40 | 6.0 |
| 1,2-Dichloroethane | | 0.55 | U | 0.55 | 6.0 |
| 1,2-Dichloropropane | | 0.17 | U | 0.17 | 6.0 |
| 1,3-Dichlorobenzene | | 0.82 | U | 0.82 | 6.0 |
| 1,3-Dichloropropane | | 0.48 | U | 0.48 | 6.0 |
| 1,4-Dichlorobenzene | | 0.33 | U | 0.33 | 6.0 |
| 2,2-Dichloropropane | | 0.52 | U | 0.52 | 6.0 |
| 2-Butanone (MEK) | | 3.9 | J B | 1.3 | 24 |
| 2-Chloroethyl vinyl ether | | 0.76 | U | 0.76 | 60 |
| 2-Chlorotoluene | | 0.36 | U | 0.36 | 6.0 |
| 2-Hexanone | | 0.81 | U | 0.81 | 24 |
| 4-Chlorotoluene | | 0.31 | U | 0.31 | 6.0 |
| p-Isopropyltoluene | | 0.49 | U | 0.49 | 6.0 |
| 4-Methyl-2-pentanone (MIBK) | | 1.5 | U | 1.5 | 24 |
| Acetone | | 6.8 | U | 6.8 | 24 |
| Benzene | | 1.1 | U | 1.1 | 6.0 |
| Bromobenzene | | 0.39 | U | 0.39 | 6.0 |
| Bromoform | | 0.29 | U | 0.29 | 6.0 |
| Bromomethane | | 0.50 | U | 0.50 | 6.0 |
| Carbon disulfide | | 0.70 | U | 0.70 | 6.0 |
| Carbon tetrachloride | | 0.77 | U | 0.77 | 6.0 |
| Chlorobenzene | | 0.63 | U | 0.63 | 6.0 |
| Dibromochloromethane | | 0.44 | U | 0.44 | 6.0 |
| Chloroethane | | 0.52 | U | 0.52 | 6.0 |
| Chloroform | | 0.43 | U | 0.43 | 6.0 |
| Chloromethane | | 0.94 | U | 0.94 | 6.0 |
| cis-1,2-Dichloroethene | | 0.33 | U | 0.33 | 6.0 |
| cis-1,3-Dichloropropene | | 0.85 | U | 0.85 | 6.0 |
| Dibromomethane | | 0.50 | U | 0.50 | 6.0 |
| Bromochloromethane | | 0.39 | U | 0.39 | 6.0 |
| Dichlorodifluoromethane | | 0.33 | U | 0.33 | 6.0 |
| Bromodichloromethane | | 0.35 | U | 0.35 | 6.0 |
| tert-Butyl alcohol (TBA) | | 6.8 | U | 6.8 | 240 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-11-14**

Lab Sample ID: 240-57597-11 Date Sampled: 11/05/2015 0920
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146348.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.03 g |
| Analysis Date: | 11/10/2015 1901 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.38 | U | 0.38 | 6.0 |
| Hexachlorobutadiene | | 0.48 | U | 0.48 | 6.0 |
| Diisopropyl ether | | 0.70 | U | 0.70 | 12 |
| Isopropylbenzene | | 0.24 | U | 0.24 | 6.0 |
| Methyl tert-butyl ether | | 0.51 | U | 0.51 | 6.0 |
| Methylene Chloride | | 0.88 | U | 0.88 | 6.0 |
| m-Xylene & p-Xylene | | 0.50 | U | 0.50 | 12 |
| Naphthalene | | 0.42 | U | 0.42 | 6.0 |
| n-Butylbenzene | | 0.70 | U | 0.70 | 6.0 |
| N-Propylbenzene | | 0.71 | U | 0.71 | 6.0 |
| o-Xylene | | 0.68 | U | 0.68 | 6.0 |
| sec-Butylbenzene | | 1.1 | U | 1.1 | 6.0 |
| Styrene | | 0.48 | U | 0.48 | 6.0 |
| Tert-amyl methyl ether | | 0.51 | U | 0.51 | 6.0 |
| Ethyl t-butyl ether | | 0.49 | U | 0.49 | 6.0 |
| Tetrachloroethene | | 0.95 | U | 0.95 | 6.0 |
| Toluene | | 0.32 | U | 0.32 | 6.0 |
| trans-1,2-Dichloroethene | | 0.38 | U | 0.38 | 6.0 |
| trans-1,3-Dichloropropene | | 0.35 | U | 0.35 | 6.0 |
| Trichloroethene | | 0.45 | U | 0.45 | 6.0 |
| Trichlorofluoromethane | | 0.35 | U | 0.35 | 6.0 |
| Vinyl acetate | | 0.79 | U | 0.79 | 12 |
| Vinyl chloride | | 0.36 | U | 0.36 | 6.0 |
| Xylenes, Total | | 0.64 | U | 0.64 | 12 |
| tert-Butylbenzene | | 0.80 | U | 0.80 | 6.0 |
| 1,2-Dibromoethane | | 0.63 | U | 0.63 | 6.0 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 90 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 87 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 81 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 91 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-12-14**

Lab Sample ID: 240-57597-12 Date Sampled: 11/05/2015 0922
Client Matrix: Solid % Moisture: 15.9 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146349.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.37 g |
| Analysis Date: | 11/10/2015 1922 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.55 | U | 0.55 | 5.5 |
| 1,1,1-Trichloroethane | | 0.72 | U | 0.72 | 5.5 |
| 1,1,2,2-Tetrachloroethane | | 0.34 | U | 0.34 | 5.5 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.46 | U | 0.46 | 5.5 |
| 1,1-Dichloroethane | | 0.38 | U | 0.38 | 5.5 |
| 1,1-Dichloroethene | | 0.89 | U | 0.89 | 5.5 |
| 1,1-Dichloropropene | | 0.41 | U | 0.41 | 5.5 |
| 1,2,3-Trichlorobenzene | | 0.55 | U | 0.55 | 5.5 |
| 1,2,3-Trichloropropane | | 0.86 | U | 0.86 | 5.5 |
| 1,2,3-Trimethylbenzene | | 0.51 | U | 0.51 | 5.5 |
| 1,2,4-Trichlorobenzene | | 0.42 | U | 0.42 | 5.5 |
| 1,2,4-Trimethylbenzene | | 0.64 | U | 0.64 | 5.5 |
| 1,2-Dibromo-3-Chloropropane | | 3.0 | U | 3.0 | 11 |
| 1,2-Dichlorobenzene | | 0.38 | U | 0.38 | 5.5 |
| 1,2-Dichloroethane | | 0.51 | U | 0.51 | 5.5 |
| 1,2-Dichloropropane | | 0.15 | U | 0.15 | 5.5 |
| 1,3-Dichlorobenzene | | 0.76 | U | 0.76 | 5.5 |
| 1,3-Dichloropropane | | 0.44 | U | 0.44 | 5.5 |
| 1,4-Dichlorobenzene | | 0.31 | U | 0.31 | 5.5 |
| 2,2-Dichloropropane | | 0.49 | U | 0.49 | 5.5 |
| 2-Butanone (MEK) | | 3.7 | J B | 1.2 | 22 |
| 2-Chloroethyl vinyl ether | | 0.71 | U | 0.71 | 55 |
| 2-Chlorotoluene | | 0.33 | U | 0.33 | 5.5 |
| 2-Hexanone | | 0.75 | U | 0.75 | 22 |
| 4-Chlorotoluene | | 0.29 | U | 0.29 | 5.5 |
| p-Isopropyltoluene | | 0.45 | U | 0.45 | 5.5 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 22 |
| Acetone | | 6.3 | U | 6.3 | 22 |
| Benzene | | 1.0 | U | 1.0 | 5.5 |
| Bromobenzene | | 0.37 | U | 0.37 | 5.5 |
| Bromoform | | 0.27 | U | 0.27 | 5.5 |
| Bromomethane | | 0.46 | U | 0.46 | 5.5 |
| Carbon disulfide | | 0.65 | U | 0.65 | 5.5 |
| Carbon tetrachloride | | 0.72 | U | 0.72 | 5.5 |
| Chlorobenzene | | 0.59 | U | 0.59 | 5.5 |
| Dibromochloromethane | | 0.41 | U | 0.41 | 5.5 |
| Chloroethane | | 0.49 | U | 0.49 | 5.5 |
| Chloroform | | 0.40 | U | 0.40 | 5.5 |
| Chloromethane | | 0.87 | U | 0.87 | 5.5 |
| cis-1,2-Dichloroethene | | 0.31 | U | 0.31 | 5.5 |
| cis-1,3-Dichloropropene | | 0.79 | U | 0.79 | 5.5 |
| Dibromomethane | | 0.46 | U | 0.46 | 5.5 |
| Bromochloromethane | | 0.37 | U | 0.37 | 5.5 |
| Dichlorodifluoromethane | | 0.31 | U | 0.31 | 5.5 |
| Bromodichloromethane | | 0.32 | U | 0.32 | 5.5 |
| tert-Butyl alcohol (TBA) | | 6.3 | U | 6.3 | 220 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-12-14**

Lab Sample ID: 240-57597-12 Date Sampled: 11/05/2015 0922
Client Matrix: Solid % Moisture: 15.9 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146349.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.37 g |
| Analysis Date: | 11/10/2015 1922 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.35 | U | 0.35 | 5.5 |
| Hexachlorobutadiene | | 0.44 | U | 0.44 | 5.5 |
| Diisopropyl ether | | 0.65 | U | 0.65 | 11 |
| Isopropylbenzene | | 0.22 | U | 0.22 | 5.5 |
| Methyl tert-butyl ether | | 0.48 | U | 0.48 | 5.5 |
| Methylene Chloride | | 0.82 | U | 0.82 | 5.5 |
| m-Xylene & p-Xylene | | 0.46 | U | 0.46 | 11 |
| Naphthalene | | 0.39 | U | 0.39 | 5.5 |
| n-Butylbenzene | | 0.65 | U | 0.65 | 5.5 |
| N-Propylbenzene | | 0.66 | U | 0.66 | 5.5 |
| o-Xylene | | 0.63 | U | 0.63 | 5.5 |
| sec-Butylbenzene | | 1.0 | U | 1.0 | 5.5 |
| Styrene | | 0.44 | U | 0.44 | 5.5 |
| Tert-amyl methyl ether | | 0.48 | U | 0.48 | 5.5 |
| Ethyl t-butyl ether | | 0.45 | U | 0.45 | 5.5 |
| Tetrachloroethene | | 0.89 | U | 0.89 | 5.5 |
| Toluene | | 0.30 | J | 0.30 | 5.5 |
| trans-1,2-Dichloroethene | | 0.35 | U | 0.35 | 5.5 |
| trans-1,3-Dichloropropene | | 0.32 | U | 0.32 | 5.5 |
| Trichloroethene | | 0.42 | U | 0.42 | 5.5 |
| Trichlorofluoromethane | | 0.32 | U | 0.32 | 5.5 |
| Vinyl acetate | | 0.73 | U | 0.73 | 11 |
| Vinyl chloride | | 0.33 | U | 0.33 | 5.5 |
| Xylenes, Total | | 0.60 | U | 0.60 | 11 |
| tert-Butylbenzene | | 0.74 | U | 0.74 | 5.5 |
| 1,2-Dibromoethane | | 0.59 | U | 0.59 | 5.5 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 99 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 98 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 87 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 103 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-13-14**

Lab Sample ID: 240-57597-13 Date Sampled: 11/05/2015 0924
Client Matrix: Solid % Moisture: 14.7 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146350.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.67 g |
| Analysis Date: | 11/10/2015 1944 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.52 | U | 0.52 | 5.2 |
| 1,1,1-Trichloroethane | | 0.67 | U | 0.67 | 5.2 |
| 1,1,2,2-Tetrachloroethane | | 0.32 | U | 0.32 | 5.2 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.43 | U | 0.43 | 5.2 |
| 1,1-Dichloroethane | | 0.35 | U | 0.35 | 5.2 |
| 1,1-Dichloroethene | | 0.83 | U | 0.83 | 5.2 |
| 1,1-Dichloropropene | | 0.38 | U | 0.38 | 5.2 |
| 1,2,3-Trichlorobenzene | | 0.52 | U | 0.52 | 5.2 |
| 1,2,3-Trichloropropane | | 0.81 | U | 0.81 | 5.2 |
| 1,2,3-Trimethylbenzene | | 0.48 | U | 0.48 | 5.2 |
| 1,2,4-Trichlorobenzene | | 0.39 | U | 0.39 | 5.2 |
| 1,2,4-Trimethylbenzene | | 0.60 | U | 0.60 | 5.2 |
| 1,2-Dibromo-3-Chloropropane | | 2.8 | U | 2.8 | 10 |
| 1,2-Dichlorobenzene | | 0.35 | U | 0.35 | 5.2 |
| 1,2-Dichloroethane | | 0.48 | U | 0.48 | 5.2 |
| 1,2-Dichloropropane | | 0.14 | U | 0.14 | 5.2 |
| 1,3-Dichlorobenzene | | 0.71 | U | 0.71 | 5.2 |
| 1,3-Dichloropropane | | 0.41 | U | 0.41 | 5.2 |
| 1,4-Dichlorobenzene | | 0.29 | U | 0.29 | 5.2 |
| 2,2-Dichloropropane | | 0.45 | U | 0.45 | 5.2 |
| 2-Butanone (MEK) | | 3.1 | J B | 1.2 | 21 |
| 2-Chloroethyl vinyl ether | | 0.66 | U | 0.66 | 52 |
| 2-Chlorotoluene | | 0.31 | U | 0.31 | 5.2 |
| 2-Hexanone | | 0.70 | U | 0.70 | 21 |
| 4-Chlorotoluene | | 0.27 | U | 0.27 | 5.2 |
| p-Isopropyltoluene | | 0.42 | U | 0.42 | 5.2 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 21 |
| Acetone | | 5.9 | U | 5.9 | 21 |
| Benzene | | 0.93 | U | 0.93 | 5.2 |
| Bromobenzene | | 0.34 | U | 0.34 | 5.2 |
| Bromoform | | 0.25 | U | 0.25 | 5.2 |
| Bromomethane | | 0.43 | U | 0.43 | 5.2 |
| Carbon disulfide | | 0.61 | U | 0.61 | 5.2 |
| Carbon tetrachloride | | 0.67 | U | 0.67 | 5.2 |
| Chlorobenzene | | 0.55 | U | 0.55 | 5.2 |
| Dibromochloromethane | | 0.38 | U | 0.38 | 5.2 |
| Chloroethane | | 0.45 | U | 0.45 | 5.2 |
| Chloroform | | 0.37 | U | 0.37 | 5.2 |
| Chloromethane | | 0.82 | U | 0.82 | 5.2 |
| cis-1,2-Dichloroethene | | 0.29 | U | 0.29 | 5.2 |
| cis-1,3-Dichloropropene | | 0.73 | U | 0.73 | 5.2 |
| Dibromomethane | | 0.43 | U | 0.43 | 5.2 |
| Bromochloromethane | | 0.34 | U | 0.34 | 5.2 |
| Dichlorodifluoromethane | | 0.29 | U | 0.29 | 5.2 |
| Bromodichloromethane | | 0.30 | U | 0.30 | 5.2 |
| tert-Butyl alcohol (TBA) | | 5.9 | U | 5.9 | 210 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-13-14**

Lab Sample ID: 240-57597-13 Date Sampled: 11/05/2015 0924
Client Matrix: Solid % Moisture: 14.7 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146350.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.67 g |
| Analysis Date: | 11/10/2015 1944 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.33 | U | 0.33 | 5.2 |
| Hexachlorobutadiene | | 0.41 | U | 0.41 | 5.2 |
| Diisopropyl ether | | 0.61 | U | 0.61 | 10 |
| Isopropylbenzene | | 0.21 | U | 0.21 | 5.2 |
| Methyl tert-butyl ether | | 0.44 | U | 0.44 | 5.2 |
| Methylene Chloride | | 0.76 | U | 0.76 | 5.2 |
| m-Xylene & p-Xylene | | 0.43 | U | 0.43 | 10 |
| Naphthalene | | 0.36 | U | 0.36 | 5.2 |
| n-Butylbenzene | | 0.61 | U | 0.61 | 5.2 |
| N-Propylbenzene | | 0.62 | U | 0.62 | 5.2 |
| o-Xylene | | 0.59 | U | 0.59 | 5.2 |
| sec-Butylbenzene | | 0.94 | U | 0.94 | 5.2 |
| Styrene | | 0.41 | U | 0.41 | 5.2 |
| Tert-amyl methyl ether | | 0.44 | U | 0.44 | 5.2 |
| Ethyl t-butyl ether | | 0.42 | U | 0.42 | 5.2 |
| Tetrachloroethene | | 0.83 | U | 0.83 | 5.2 |
| Toluene | | 0.28 | U | 0.28 | 5.2 |
| trans-1,2-Dichloroethene | | 0.33 | U | 0.33 | 5.2 |
| trans-1,3-Dichloropropene | | 0.30 | U | 0.30 | 5.2 |
| Trichloroethene | | 0.39 | U | 0.39 | 5.2 |
| Trichlorofluoromethane | | 0.30 | U | 0.30 | 5.2 |
| Vinyl acetate | | 0.68 | U | 0.68 | 10 |
| Vinyl chloride | | 0.31 | U | 0.31 | 5.2 |
| Xylenes, Total | | 0.56 | U | 0.56 | 10 |
| tert-Butylbenzene | | 0.69 | U | 0.69 | 5.2 |
| 1,2-Dibromoethane | | 0.55 | U | 0.55 | 5.2 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 105 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 98 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 90 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 100 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-14-14**

Lab Sample ID: 240-57597-14 Date Sampled: 11/05/2015 0926
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146351.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/10/2015 2005 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.53 | U | 0.53 | 5.3 |
| 1,1,1-Trichloroethane | | 0.68 | U | 0.68 | 5.3 |
| 1,1,2,2-Tetrachloroethane | | 0.33 | U | 0.33 | 5.3 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.44 | U | 0.44 | 5.3 |
| 1,1-Dichloroethane | | 0.36 | U | 0.36 | 5.3 |
| 1,1-Dichloroethene | | 0.84 | U | 0.84 | 5.3 |
| 1,1-Dichloropropene | | 0.39 | U | 0.39 | 5.3 |
| 1,2,3-Trichlorobenzene | | 0.53 | U | 0.53 | 5.3 |
| 1,2,3-Trichloropropane | | 0.82 | U | 0.82 | 5.3 |
| 1,2,3-Trimethylbenzene | | 0.48 | U | 0.48 | 5.3 |
| 1,2,4-Trichlorobenzene | | 0.40 | U | 0.40 | 5.3 |
| 1,2,4-Trimethylbenzene | | 0.61 | U | 0.61 | 5.3 |
| 1,2-Dibromo-3-Chloropropane | | 2.9 | U | 2.9 | 11 |
| 1,2-Dichlorobenzene | | 0.36 | U | 0.36 | 5.3 |
| 1,2-Dichloroethane | | 0.48 | U | 0.48 | 5.3 |
| 1,2-Dichloropropane | | 0.15 | U | 0.15 | 5.3 |
| 1,3-Dichlorobenzene | | 0.73 | U | 0.73 | 5.3 |
| 1,3-Dichloropropane | | 0.42 | U | 0.42 | 5.3 |
| 1,4-Dichlorobenzene | | 0.30 | U | 0.30 | 5.3 |
| 2,2-Dichloropropane | | 0.46 | U | 0.46 | 5.3 |
| 2-Butanone (MEK) | | 4.0 | J B | 1.2 | 21 |
| 2-Chloroethyl vinyl ether | | 0.67 | U | 0.67 | 53 |
| 2-Chlorotoluene | | 0.32 | U | 0.32 | 5.3 |
| 2-Hexanone | | 0.72 | U | 0.72 | 21 |
| 4-Chlorotoluene | | 0.27 | U | 0.27 | 5.3 |
| p-Isopropyltoluene | | 0.43 | U | 0.43 | 5.3 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 21 |
| Acetone | | 6.0 | U | 6.0 | 21 |
| Benzene | | 0.95 | U | 0.95 | 5.3 |
| Bromobenzene | | 0.35 | U | 0.35 | 5.3 |
| Bromoform | | 0.25 | U | 0.25 | 5.3 |
| Bromomethane | | 0.44 | U | 0.44 | 5.3 |
| Carbon disulfide | | 0.62 | U | 0.62 | 5.3 |
| Carbon tetrachloride | | 0.68 | U | 0.68 | 5.3 |
| Chlorobenzene | | 0.56 | U | 0.56 | 5.3 |
| Dibromochloromethane | | 0.39 | U | 0.39 | 5.3 |
| Chloroethane | | 0.46 | U | 0.46 | 5.3 |
| Chloroform | | 0.38 | U | 0.38 | 5.3 |
| Chloromethane | | 0.83 | U | 0.83 | 5.3 |
| cis-1,2-Dichloroethene | | 0.30 | U | 0.30 | 5.3 |
| cis-1,3-Dichloropropene | | 0.75 | U | 0.75 | 5.3 |
| Dibromomethane | | 0.44 | U | 0.44 | 5.3 |
| Bromochloromethane | | 0.35 | U | 0.35 | 5.3 |
| Dichlorodifluoromethane | | 0.30 | U | 0.30 | 5.3 |
| Bromodichloromethane | | 0.31 | U | 0.31 | 5.3 |
| tert-Butyl alcohol (TBA) | | 6.0 | U | 6.0 | 210 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-14-14**

Lab Sample ID: 240-57597-14 Date Sampled: 11/05/2015 0926
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206089 | Lab File ID: | 146351.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/10/2015 2005 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.34 | U | 0.34 | 5.3 |
| Hexachlorobutadiene | | 0.42 | U | 0.42 | 5.3 |
| Diisopropyl ether | | 0.62 | U | 0.62 | 11 |
| Isopropylbenzene | | 0.21 | U | 0.21 | 5.3 |
| Methyl tert-butyl ether | | 0.45 | U | 0.45 | 5.3 |
| Methylene Chloride | | 0.78 | U | 0.78 | 5.3 |
| m-Xylene & p-Xylene | | 0.44 | U | 0.44 | 11 |
| Naphthalene | | 0.37 | U | 0.37 | 5.3 |
| n-Butylbenzene | | 0.62 | U | 0.62 | 5.3 |
| N-Propylbenzene | | 0.63 | U | 0.63 | 5.3 |
| o-Xylene | | 0.60 | U | 0.60 | 5.3 |
| sec-Butylbenzene | | 0.96 | U | 0.96 | 5.3 |
| Styrene | | 0.42 | U | 0.42 | 5.3 |
| Tert-amyl methyl ether | | 0.45 | U | 0.45 | 5.3 |
| Ethyl t-butyl ether | | 0.43 | U | 0.43 | 5.3 |
| Tetrachloroethene | | 0.84 | U | 0.84 | 5.3 |
| Toluene | | 0.29 | J | 0.28 | 5.3 |
| trans-1,2-Dichloroethene | | 0.34 | U | 0.34 | 5.3 |
| trans-1,3-Dichloropropene | | 0.31 | U | 0.31 | 5.3 |
| Trichloroethene | | 0.40 | U | 0.40 | 5.3 |
| Trichlorofluoromethane | | 0.31 | U | 0.31 | 5.3 |
| Vinyl acetate | | 0.70 | U | 0.70 | 11 |
| Vinyl chloride | | 0.32 | U | 0.32 | 5.3 |
| Xylenes, Total | | 0.57 | U | 0.57 | 11 |
| tert-Butylbenzene | | 0.71 | U | 0.71 | 5.3 |
| 1,2-Dibromoethane | | 0.56 | U | 0.56 | 5.3 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 87 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 89 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 78 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 96 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-DUP**

Lab Sample ID: 240-57597-15 Date Sampled: 11/05/2015 0000
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206305 | Lab File ID: | 146365.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/11/2015 1309 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 1103 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.53 | U | 0.53 | 5.3 |
| 1,1,1-Trichloroethane | | 0.68 | U | 0.68 | 5.3 |
| 1,1,2,2-Tetrachloroethane | | 0.33 | U | 0.33 | 5.3 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.44 | U | 0.44 | 5.3 |
| 1,1-Dichloroethane | | 0.36 | U | 0.36 | 5.3 |
| 1,1-Dichloroethene | | 0.84 | U | 0.84 | 5.3 |
| 1,1-Dichloropropene | | 0.39 | U | 0.39 | 5.3 |
| 1,2,3-Trichlorobenzene | | 0.53 | U | 0.53 | 5.3 |
| 1,2,3-Trichloropropane | | 0.82 | U | 0.82 | 5.3 |
| 1,2,3-Trimethylbenzene | | 0.48 | U | 0.48 | 5.3 |
| 1,2,4-Trichlorobenzene | | 0.40 | U | 0.40 | 5.3 |
| 1,2,4-Trimethylbenzene | | 0.61 | U | 0.61 | 5.3 |
| 1,2-Dibromo-3-Chloropropane | | 2.9 | U | 2.9 | 11 |
| 1,2-Dichlorobenzene | | 0.36 | U | 0.36 | 5.3 |
| 1,2-Dichloroethane | | 0.48 | U | 0.48 | 5.3 |
| 1,2-Dichloropropane | | 0.15 | U | 0.15 | 5.3 |
| 1,3-Dichlorobenzene | | 0.73 | U | 0.73 | 5.3 |
| 1,3-Dichloropropane | | 0.42 | U | 0.42 | 5.3 |
| 1,4-Dichlorobenzene | | 0.29 | U | 0.29 | 5.3 |
| 2,2-Dichloropropane | | 0.46 | U | 0.46 | 5.3 |
| 2-Butanone (MEK) | | 3.6 | J B | 1.2 | 21 |
| 2-Chloroethyl vinyl ether | | 0.67 | U | 0.67 | 53 |
| 2-Chlorotoluene | | 0.32 | U | 0.32 | 5.3 |
| 2-Hexanone | | 0.72 | U | 0.72 | 21 |
| 4-Chlorotoluene | | 0.27 | U | 0.27 | 5.3 |
| p-Isopropyltoluene | | 0.43 | U | 0.43 | 5.3 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 21 |
| Acetone | | 6.0 | U | 6.0 | 21 |
| Benzene | | 0.95 | U | 0.95 | 5.3 |
| Bromobenzene | | 0.35 | U | 0.35 | 5.3 |
| Bromoform | | 0.25 | U | 0.25 | 5.3 |
| Bromomethane | | 0.44 | U | 0.44 | 5.3 |
| Carbon disulfide | | 0.62 | U | 0.62 | 5.3 |
| Carbon tetrachloride | | 0.68 | U | 0.68 | 5.3 |
| Chlorobenzene | | 0.56 | U | 0.56 | 5.3 |
| Dibromochloromethane | | 0.39 | U | 0.39 | 5.3 |
| Chloroethane | | 0.46 | U | 0.46 | 5.3 |
| Chloroform | | 0.38 | U | 0.38 | 5.3 |
| Chloromethane | | 0.83 | U | 0.83 | 5.3 |
| cis-1,2-Dichloroethene | | 0.29 | U | 0.29 | 5.3 |
| cis-1,3-Dichloropropene | | 0.75 | U | 0.75 | 5.3 |
| Dibromomethane | | 0.44 | U | 0.44 | 5.3 |
| Bromochloromethane | | 0.35 | U | 0.35 | 5.3 |
| Dichlorodifluoromethane | | 0.29 | U | 0.29 | 5.3 |
| Bromodichloromethane | | 0.31 | U | 0.31 | 5.3 |
| tert-Butyl alcohol (TBA) | | 6.0 | U | 6.0 | 210 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-DUP**

Lab Sample ID: 240-57597-15 Date Sampled: 11/05/2015 0000
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-206305 | Lab File ID: | 146365.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/11/2015 1309 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 1103 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.34 | U | 0.34 | 5.3 |
| Hexachlorobutadiene | | 0.42 | U | 0.42 | 5.3 |
| Diisopropyl ether | | 0.62 | U | 0.62 | 11 |
| Isopropylbenzene | | 0.21 | U | 0.21 | 5.3 |
| Methyl tert-butyl ether | | 0.45 | U | 0.45 | 5.3 |
| Methylene Chloride | | 0.78 | U | 0.78 | 5.3 |
| m-Xylene & p-Xylene | | 0.44 | U | 0.44 | 11 |
| Naphthalene | | 0.37 | U | 0.37 | 5.3 |
| n-Butylbenzene | | 0.62 | U | 0.62 | 5.3 |
| N-Propylbenzene | | 0.63 | U | 0.63 | 5.3 |
| o-Xylene | | 0.60 | U | 0.60 | 5.3 |
| sec-Butylbenzene | | 0.96 | U | 0.96 | 5.3 |
| Styrene | | 0.42 | U | 0.42 | 5.3 |
| Tert-amyl methyl ether | | 0.45 | U | 0.45 | 5.3 |
| Ethyl t-butyl ether | | 0.43 | U | 0.43 | 5.3 |
| Tetrachloroethene | | 0.84 | U | 0.84 | 5.3 |
| Toluene | | 0.28 | U | 0.28 | 5.3 |
| trans-1,2-Dichloroethene | | 0.34 | U | 0.34 | 5.3 |
| trans-1,3-Dichloropropene | | 0.31 | U | 0.31 | 5.3 |
| Trichloroethene | | 0.40 | U | 0.40 | 5.3 |
| Trichlorofluoromethane | | 0.31 | U | 0.31 | 5.3 |
| Vinyl acetate | | 0.70 | U | 0.70 | 11 |
| Vinyl chloride | | 0.32 | U | 0.32 | 5.3 |
| Xylenes, Total | | 0.57 | U | 0.57 | 11 |
| tert-Butylbenzene | | 0.71 | U | 0.71 | 5.3 |
| 1,2-Dibromoethane | | 0.56 | U | 0.56 | 5.3 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 102 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 101 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 89 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 102 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **TB-110515**

Lab Sample ID: 240-57597-16 Date Sampled: 11/05/2015 0000
Client Matrix: Water Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Prep Method: | 5030C | Prep Batch: | N/A | Lab File ID: | UXC4397.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 1233 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 1233 | | | | |

| Analyte | Result (ug/L) | Qualifier | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.28 | U | 0.28 | 1.0 |
| 1,1,1-Trichloroethane | 0.44 | U | 0.44 | 1.0 |
| 1,1,2,2-Tetrachloroethane | 0.22 | U | 0.22 | 1.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloroethane | 0.30 | U | 0.30 | 1.0 |
| 1,1-Dichloroethene | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloropropene | 0.42 | U | 0.42 | 1.0 |
| 1,2,3-Trichlorobenzene | 0.37 | U | 0.37 | 1.0 |
| 1,2,3-Trichloropropane | 0.44 | U | 0.44 | 1.0 |
| 1,2,3-Trimethylbenzene | 0.47 | U | 0.47 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.32 | U | 0.32 | 1.0 |
| 1,2,4-Trimethylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromo-3-Chloropropane | 0.82 | U | 0.82 | 2.0 |
| 1,2-Dichlorobenzene | 0.25 | U | 0.25 | 1.0 |
| 1,2-Dichloroethane | 0.23 | U | 0.23 | 1.0 |
| 1,2-Dichloropropane | 0.25 | U | 0.25 | 1.0 |
| 1,3-Dichlorobenzene | 0.19 | U | 0.19 | 1.0 |
| 1,3-Dichloropropane | 0.19 | U | 0.19 | 1.0 |
| 1,4-Dichlorobenzene | 0.27 | U | 0.27 | 1.0 |
| 2,2-Dichloropropane | 0.26 | U | 0.26 | 1.0 |
| 2-Butanone (MEK) | 0.53 | U | 0.53 | 10 |
| 2-Chloroethyl vinyl ether | 0.59 | U | 0.59 | 10 |
| 2-Chlorotoluene | 0.40 | U | 0.40 | 1.0 |
| 2-Hexanone | 0.48 | U | 0.48 | 10 |
| 4-Chlorotoluene | 0.29 | U | 0.29 | 1.0 |
| p-Isopropyltoluene | 0.43 | U | 0.43 | 1.0 |
| 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 0.99 | 10 |
| Acetone | 0.94 | U | 0.94 | 10 |
| Benzene | 0.35 | U | 0.35 | 1.0 |
| Bromobenzene | 0.35 | U | 0.35 | 1.0 |
| Bromoform | 0.56 | U | 0.56 | 1.0 |
| Bromomethane | 0.44 | U | 0.44 | 1.0 |
| Carbon disulfide | 0.38 | U | 0.38 | 1.0 |
| Carbon tetrachloride | 0.43 | U | 0.43 | 1.0 |
| Chlorobenzene | 0.25 | U | 0.25 | 1.0 |
| Dibromochloromethane | 0.43 | U | 0.43 | 1.0 |
| Chloroethane | 0.32 | U | 0.32 | 1.0 |
| Chloroform | 0.25 | U | 0.25 | 1.0 |
| Chloromethane | 0.44 | U | 0.44 | 1.0 |
| cis-1,2-Dichloroethene | 0.26 | U | 0.26 | 1.0 |
| cis-1,3-Dichloropropene | 0.46 | U | 0.46 | 1.0 |
| Dibromomethane | 0.42 | U | 0.42 | 1.0 |
| Bromochloromethane | 0.50 | U | 0.50 | 1.0 |
| Dichlorodifluoromethane | 0.32 | U | 0.32 | 1.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 1.0 |
| tert-Butyl alcohol (TBA) | 4.9 | U | 4.9 | 50 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: TB-110515**

Lab Sample ID: 240-57597-16 Date Sampled: 11/05/2015 0000
Client Matrix: Water Date Received: 11/06/2015 1020

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Prep Method: | 5030C | Prep Batch: | N/A | Lab File ID: | UXC4397.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 1233 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 1233 | | | | |

| Analyte | Result (ug/L) | Qualifier | MDL | RL |
|---|---|---|---|---|
| Ethylbenzene | 0.25 | U | 0.25 | 1.0 |
| Hexachlorobutadiene | 0.35 | U | 0.35 | 1.0 |
| Diisopropyl ether | 0.50 | U | 0.50 | 10 |
| Isopropylbenzene | 0.35 | U | 0.35 | 1.0 |
| Methyl tert-butyl ether | 0.20 | U | 0.20 | 1.0 |
| Methylene Chloride | 0.33 | U | 0.33 | 1.0 |
| m-Xylene & p-Xylene | 0.24 | U | 0.24 | 2.0 |
| Naphthalene | 0.45 | U | 0.45 | 1.0 |
| n-Butylbenzene | 0.31 | U | 0.31 | 1.0 |
| N-Propylbenzene | 0.40 | U | 0.40 | 1.0 |
| o-Xylene | 0.25 | U | 0.25 | 1.0 |
| sec-Butylbenzene | 0.48 | U | 0.48 | 1.0 |
| Styrene | 0.45 | U | 0.45 | 1.0 |
| Tert-amyl methyl ether | 0.30 | U | 0.30 | 5.0 |
| Ethyl t-butyl ether | 0.23 | U | 0.23 | 5.0 |
| Tetrachloroethene | 0.31 | U | 0.31 | 1.0 |
| Toluene | 0.23 | U | 0.23 | 1.0 |
| trans-1,2-Dichloroethene | 0.30 | U | 0.30 | 1.0 |
| trans-1,3-Dichloropropene | 0.56 | U | 0.56 | 1.0 |
| Trichloroethene | 0.22 | U | 0.22 | 1.0 |
| Trichlorofluoromethane | 0.49 | U | 0.49 | 1.0 |
| Vinyl acetate | 0.41 | U | 0.41 | 2.0 |
| Vinyl chloride | 0.29 | U | 0.29 | 1.0 |
| Xylenes, Total | 0.52 | U | 0.52 | 2.0 |
| tert-Butylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromoethane | 0.32 | U | 0.32 | 1.0 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 85 | | 80 - 120 |
| Dibromofluoromethane (Surr) | 97 | | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 90 | | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 99 | | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-1-14**

Lab Sample ID: 240-57597-1 Date Sampled: 11/05/2015 0900
Client Matrix: Solid % Moisture: 15.7 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111006.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.99 g |
| Analysis Date: | 11/11/2015 1116 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 59 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.53 | U | 0.53 | 59 |
| 2-Chlorophenol | | 9.7 | U | 9.7 | 59 |
| 2-Methylnaphthalene | | 0.59 | U | 0.59 | 7.9 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.9 | U | 9.9 | 59 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.90 | U | 0.90 | 7.9 |
| Acenaphthylene | | 0.42 | U | 0.42 | 7.9 |
| Acetophenone | | 16 | J | 11 | 120 |
| Anthracene | | 0.93 | U | 0.93 | 7.9 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.75 | U | 0.75 | 7.9 |
| Benzo[a]pyrene | | 0.76 | U | 0.76 | 7.9 |
| Benzo[b]fluoranthene | | 0.70 | U | 0.70 | 7.9 |
| Benzo[g,h,i]perylene | | 0.42 | U | 0.42 | 7.9 |
| Benzo[k]fluoranthene | | 0.81 | U | 0.81 | 7.9 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 31 | J | 23 | 83 |
| Butyl benzyl phthalate | | 12 | U | 12 | 83 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 59 |
| Chrysene | | 1.3 | U | 1.3 | 7.9 |
| Dibenz(a,h)anthracene | | 0.78 | U | 0.78 | 7.9 |
| Dibenzofuran | | 0.78 | U | 0.78 | 59 |
| Diethyl phthalate | | 19 | U | 19 | 83 |
| Dimethyl phthalate | | 20 | U | 20 | 83 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-1-14**

Lab Sample ID: 240-57597-1 Date Sampled: 11/05/2015 0900
Client Matrix: Solid % Moisture: 15.7 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111006.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.99 g |
| Analysis Date: | 11/11/2015 1116 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 83 |
| Di-n-octyl phthalate | | 9.4 | U | 9.4 | 83 |
| Fluoranthene | | 0.65 | U | 0.65 | 7.9 |
| Fluorene | | 0.63 | U | 0.63 | 7.9 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.9 |
| Hexachlorobutadiene | | 6.6 | U | 6.6 | 59 |
| Hexachlorocyclopentadiene | | 9.6 | U | 9.6 | 390 |
| Hexachloroethane | | 11 | U | 11 | 59 |
| Indeno[1,2,3-cd]pyrene | | 0.42 | U | 0.42 | 7.9 |
| Isophorone | | 15 | U | 15 | 59 |
| N-Nitrosodi-n-propylamine | | 7.5 | U | 7.5 | 59 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 59 |
| Naphthalene | | 0.97 | U | 0.97 | 7.9 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.87 | U | 0.87 | 7.9 |
| Phenol | | 8.7 | U | 8.7 | 59 |
| Pyrene | | 0.52 | U | 0.52 | 7.9 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 76 | | 36 - 110 |
| Phenol-d5 (Surr) | 65 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 62 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 64 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 67 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 44 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-2-14**

Lab Sample ID: 240-57597-2 Date Sampled: 11/05/2015 0902
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111007.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.16 g |
| Analysis Date: | 11/11/2015 1140 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 58 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 29 | U | 29 | 180 |
| 2,4,6-Trichlorophenol | | 10 | U | 10 | 180 |
| 2,4-Dichlorophenol | | 23 | U | 23 | 180 |
| 2,4-Dimethylphenol | | 23 | U | 23 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 230 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 230 |
| 2-Chloronaphthalene | | 0.53 | U | 0.53 | 58 |
| 2-Chlorophenol | | 9.6 | U | 9.6 | 58 |
| 2-Methylnaphthalene | | 0.58 | U | 0.58 | 7.8 |
| 2-Methylphenol | | 13 | U | 13 | 230 |
| 2-Nitroaniline | | 11 | U | 11 | 230 |
| 2-Nitrophenol | | 9.7 | U | 9.7 | 58 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 230 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Nitroaniline | | 30 | U | 30 | 230 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.89 | U | 0.89 | 7.8 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.8 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.91 | U | 0.91 | 7.8 |
| Atrazine | | 11 | U | 11 | 230 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.74 | U | 0.74 | 7.8 |
| Benzo[a]pyrene | | 0.75 | U | 0.75 | 7.8 |
| Benzo[b]fluoranthene | | 0.69 | U | 0.69 | 7.8 |
| Benzo[g,h,i]perylene | | 0.41 | U | 0.41 | 7.8 |
| Benzo[k]fluoranthene | | 0.80 | U | 0.80 | 7.8 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.3 | U | 2.3 | 120 |
| Bis(2-ethylhexyl) phthalate | | 22 | U | 22 | 82 |
| Butyl benzyl phthalate | | 12 | U | 12 | 82 |
| Caprolactam | | 43 | U | 43 | 390 |
| Carbazole | | 32 | U | 32 | 58 |
| Chrysene | | 1.3 | U | 1.3 | 7.8 |
| Dibenz(a,h)anthracene | | 0.77 | U | 0.77 | 7.8 |
| Dibenzofuran | | 0.77 | U | 0.77 | 58 |
| Diethyl phthalate | | 19 | U | 19 | 82 |
| Dimethyl phthalate | | 20 | U | 20 | 82 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-2-14**

Lab Sample ID: 240-57597-2 Date Sampled: 11/05/2015 0902
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111007.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.16 g |
| Analysis Date: | 11/11/2015 1140 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 82 |
| Di-n-octyl phthalate | | 9.2 | U | 9.2 | 82 |
| Fluoranthene | | 0.64 | U | 0.64 | 7.8 |
| Fluorene | | 0.62 | U | 0.62 | 7.8 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.8 |
| Hexachlorobutadiene | | 6.6 | U | 6.6 | 58 |
| Hexachlorocyclopentadiene | | 9.5 | U | 9.5 | 390 |
| Hexachloroethane | | 11 | U | 11 | 58 |
| Indeno[1,2,3-cd]pyrene | | 0.41 | U | 0.41 | 7.8 |
| Isophorone | | 15 | U | 15 | 58 |
| N-Nitrosodi-n-propylamine | | 7.4 | U | 7.4 | 58 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 58 |
| Naphthalene | | 0.96 | U | 0.96 | 7.8 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.85 | U | 0.85 | 7.8 |
| Phenol | | 8.5 | U | 8.5 | 58 |
| Pyrene | | 0.51 | U | 0.51 | 7.8 |
| 3 & 4 Methylphenol | | 23 | U | 23 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 78 | | 36 - 110 |
| Phenol-d5 (Surr) | 61 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 58 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 61 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 63 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 42 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-3-14**

Lab Sample ID: 240-57597-3 Date Sampled: 11/05/2015 0904
Client Matrix: Solid % Moisture: 17.8 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111008.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.03 g |
| Analysis Date: | 11/11/2015 1205 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.3 | U | 4.3 | 61 |
| bis (2-chloroisopropyl) ether | | 12 | U | 12 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 26 | U | 26 | 400 |
| 2,4-Dinitrotoluene | | 21 | U | 21 | 240 |
| 2,6-Dinitrotoluene | | 26 | U | 26 | 240 |
| 2-Chloronaphthalene | | 0.55 | U | 0.55 | 61 |
| 2-Chlorophenol | | 10 | U | 10 | 61 |
| 2-Methylnaphthalene | | 0.61 | U | 0.61 | 8.1 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 10 | U | 10 | 61 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Chloro-3-methylphenol | | 26 | U | 26 | 180 |
| 4-Chloroaniline | | 21 | U | 21 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Nitroaniline | | 32 | U | 32 | 240 |
| 4-Nitrophenol | | 21 | U | 21 | 400 |
| Acenaphthene | | 0.92 | U | 0.92 | 8.1 |
| Acenaphthylene | | 0.43 | U | 0.43 | 8.1 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.95 | U | 0.95 | 8.1 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 15 | U | 15 | 120 |
| Benzo[a]anthracene | | 0.77 | U | 0.77 | 8.1 |
| Benzo[a]pyrene | | 0.78 | U | 0.78 | 8.1 |
| Benzo[b]fluoranthene | | 0.72 | U | 0.72 | 8.1 |
| Benzo[g,h,i]perylene | | 0.43 | U | 0.43 | 8.1 |
| Benzo[k]fluoranthene | | 0.83 | U | 0.83 | 8.1 |
| Bis(2-chloroethoxy)methane | | 27 | U | 27 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 23 | U | 23 | 85 |
| Butyl benzyl phthalate | | 12 | U | 12 | 85 |
| Caprolactam | | 45 | U | 45 | 400 |
| Carbazole | | 33 | U | 33 | 61 |
| Chrysene | | 1.3 | U | 1.3 | 8.1 |
| Dibenz(a,h)anthracene | | 0.80 | U | 0.80 | 8.1 |
| Dibenzofuran | | 0.80 | U | 0.80 | 61 |
| Diethyl phthalate | | 19 | U | 19 | 85 |
| Dimethyl phthalate | | 21 | U | 21 | 85 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-3-14**

Lab Sample ID: 240-57597-3 Date Sampled: 11/05/2015 0904
Client Matrix: Solid % Moisture: 17.8 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111008.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.03 g |
| Analysis Date: | 11/11/2015 1205 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 85 |
| Di-n-octyl phthalate | | 9.6 | U | 9.6 | 85 |
| Fluoranthene | | 0.67 | U | 0.67 | 8.1 |
| Fluorene | | 0.64 | U | 0.64 | 8.1 |
| Hexachlorobenzene | | 2.6 | U | 2.6 | 8.1 |
| Hexachlorobutadiene | | 6.8 | U | 6.8 | 61 |
| Hexachlorocyclopentadiene | | 9.8 | U | 9.8 | 400 |
| Hexachloroethane | | 11 | U | 11 | 61 |
| Indeno[1,2,3-cd]pyrene | | 0.43 | U | 0.43 | 8.1 |
| Isophorone | | 16 | U | 16 | 61 |
| N-Nitrosodi-n-propylamine | | 7.7 | U | 7.7 | 61 |
| N-Nitrosodiphenylamine | | 26 | U | 26 | 61 |
| Naphthalene | | 1.0 | U | 1.0 | 8.1 |
| Nitrobenzene | | 2.7 | U | 2.7 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.89 | U | 0.89 | 8.1 |
| Phenol | | 8.9 | U | 8.9 | 61 |
| Pyrene | | 0.53 | U | 0.53 | 8.1 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 490 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 73 | | 36 - 110 |
| Phenol-d5 (Surr) | 60 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 58 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 61 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 63 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 41 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-4-14**

Lab Sample ID: 240-57597-4 Date Sampled: 11/05/2015 0906
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111009.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.84 g |
| Analysis Date: | 11/11/2015 1230 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 60 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.54 | U | 0.54 | 60 |
| 2-Chlorophenol | | 9.8 | U | 9.8 | 60 |
| 2-Methylnaphthalene | | 0.60 | U | 0.60 | 7.9 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.9 | U | 9.9 | 60 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 60 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 60 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.91 | U | 0.91 | 7.9 |
| Acenaphthylene | | 0.42 | U | 0.42 | 7.9 |
| Acetophenone | | 13 | J | 11 | 120 |
| Anthracene | | 0.93 | U | 0.93 | 7.9 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.75 | U | 0.75 | 7.9 |
| Benzo[a]pyrene | | 0.76 | U | 0.76 | 7.9 |
| Benzo[b]fluoranthene | | 0.70 | U | 0.70 | 7.9 |
| Benzo[g,h,i]perylene | | 0.42 | U | 0.42 | 7.9 |
| Benzo[k]fluoranthene | | 0.81 | U | 0.81 | 7.9 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 33 | J | 23 | 83 |
| Butyl benzyl phthalate | | 12 | U | 12 | 83 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 60 |
| Chrysene | | 1.3 | U | 1.3 | 7.9 |
| Dibenz(a,h)anthracene | | 0.79 | U | 0.79 | 7.9 |
| Dibenzofuran | | 0.79 | U | 0.79 | 60 |
| Diethyl phthalate | | 19 | U | 19 | 83 |
| Dimethyl phthalate | | 20 | U | 20 | 83 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-4-14**

Lab Sample ID: 240-57597-4 Date Sampled: 11/05/2015 0906
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111009.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.84 g |
| Analysis Date: | 11/11/2015 1230 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 83 |
| Di-n-octyl phthalate | | 9.4 | U | 9.4 | 83 |
| Fluoranthene | | 0.66 | U | 0.66 | 7.9 |
| Fluorene | | 0.63 | U | 0.63 | 7.9 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.9 |
| Hexachlorobutadiene | | 6.7 | U | 6.7 | 60 |
| Hexachlorocyclopentadiene | | 9.7 | U | 9.7 | 390 |
| Hexachloroethane | | 11 | U | 11 | 60 |
| Indeno[1,2,3-cd]pyrene | | 0.42 | U | 0.42 | 7.9 |
| Isophorone | | 15 | U | 15 | 60 |
| N-Nitrosodi-n-propylamine | | 7.5 | U | 7.5 | 60 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 60 |
| Naphthalene | | 0.98 | U | 0.98 | 7.9 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.87 | U | 0.87 | 7.9 |
| Phenol | | 8.7 | U | 8.7 | 60 |
| Pyrene | | 0.52 | U | 0.52 | 7.9 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 85 | | 36 - 110 |
| Phenol-d5 (Surr) | 72 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 69 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 72 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 72 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 42 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-5-14**

Lab Sample ID: 240-57597-5 Date Sampled: 11/05/2015 0908
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111010.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.28 g |
| Analysis Date: | 11/11/2015 1254 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 58 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 29 | U | 29 | 180 |
| 2,4,6-Trichlorophenol | | 10 | U | 10 | 180 |
| 2,4-Dichlorophenol | | 23 | U | 23 | 180 |
| 2,4-Dimethylphenol | | 23 | U | 23 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 230 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 230 |
| 2-Chloronaphthalene | | 0.53 | U | 0.53 | 58 |
| 2-Chlorophenol | | 9.6 | U | 9.6 | 58 |
| 2-Methylnaphthalene | | 0.58 | U | 0.58 | 7.8 |
| 2-Methylphenol | | 13 | U | 13 | 230 |
| 2-Nitroaniline | | 11 | U | 11 | 230 |
| 2-Nitrophenol | | 9.7 | U | 9.7 | 58 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 230 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Nitroaniline | | 30 | U | 30 | 230 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.89 | U | 0.89 | 7.8 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.8 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.91 | U | 0.91 | 7.8 |
| Atrazine | | 11 | U | 11 | 230 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.74 | U | 0.74 | 7.8 |
| Benzo[a]pyrene | | 0.75 | U | 0.75 | 7.8 |
| Benzo[b]fluoranthene | | 0.69 | U | 0.69 | 7.8 |
| Benzo[g,h,i]perylene | | 0.41 | U | 0.41 | 7.8 |
| Benzo[k]fluoranthene | | 0.79 | U | 0.79 | 7.8 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.3 | U | 2.3 | 120 |
| Bis(2-ethylhexyl) phthalate | | 38 | J | 22 | 82 |
| Butyl benzyl phthalate | | 12 | U | 12 | 82 |
| Caprolactam | | 43 | U | 43 | 390 |
| Carbazole | | 32 | U | 32 | 58 |
| Chrysene | | 1.3 | U | 1.3 | 7.8 |
| Dibenz(a,h)anthracene | | 0.77 | U | 0.77 | 7.8 |
| Dibenzofuran | | 0.77 | U | 0.77 | 58 |
| Diethyl phthalate | | 19 | U | 19 | 82 |
| Dimethyl phthalate | | 20 | U | 20 | 82 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-5-14**

Lab Sample ID: 240-57597-5 Date Sampled: 11/05/2015 0908
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111010.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.28 g |
| Analysis Date: | 11/11/2015 1254 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 82 |
| Di-n-octyl phthalate | | 9.2 | U | 9.2 | 82 |
| Fluoranthene | | 0.64 | U | 0.64 | 7.8 |
| Fluorene | | 0.62 | U | 0.62 | 7.8 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.8 |
| Hexachlorobutadiene | | 6.5 | U | 6.5 | 58 |
| Hexachlorocyclopentadiene | | 9.5 | U | 9.5 | 390 |
| Hexachloroethane | | 11 | U | 11 | 58 |
| Indeno[1,2,3-cd]pyrene | | 0.41 | U | 0.41 | 7.8 |
| Isophorone | | 15 | U | 15 | 58 |
| N-Nitrosodi-n-propylamine | | 7.4 | U | 7.4 | 58 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 58 |
| Naphthalene | | 0.96 | U | 0.96 | 7.8 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.85 | U | 0.85 | 7.8 |
| Phenol | | 8.5 | U | 8.5 | 58 |
| Pyrene | | 0.51 | U | 0.51 | 7.8 |
| 3 & 4 Methylphenol | | 23 | U | 23 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 77 | | 36 - 110 |
| Phenol-d5 (Surr) | 67 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 64 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 68 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 69 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 25 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-6-14**

Lab Sample ID: 240-57597-6 Date Sampled: 11/05/2015 0910
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111011.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.15 g |
| Analysis Date: | 11/11/2015 1319 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 60 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.54 | U | 0.54 | 60 |
| 2-Chlorophenol | | 9.8 | U | 9.8 | 60 |
| 2-Methylnaphthalene | | 0.60 | U | 0.60 | 8.0 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.9 | U | 9.9 | 60 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 60 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 60 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.91 | U | 0.91 | 8.0 |
| Acenaphthylene | | 0.42 | U | 0.42 | 8.0 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.93 | U | 0.93 | 8.0 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.75 | U | 0.75 | 8.0 |
| Benzo[a]pyrene | | 0.76 | U | 0.76 | 8.0 |
| Benzo[b]fluoranthene | | 0.70 | U | 0.70 | 8.0 |
| Benzo[g,h,i]perylene | | 0.42 | U | 0.42 | 8.0 |
| Benzo[k]fluoranthene | | 0.81 | U | 0.81 | 8.0 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 32 | J | 23 | 83 |
| Butyl benzyl phthalate | | 12 | U | 12 | 83 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 60 |
| Chrysene | | 1.3 | U | 1.3 | 8.0 |
| Dibenz(a,h)anthracene | | 0.79 | U | 0.79 | 8.0 |
| Dibenzofuran | | 0.79 | U | 0.79 | 60 |
| Diethyl phthalate | | 19 | U | 19 | 83 |
| Dimethyl phthalate | | 20 | U | 20 | 83 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-6-14**

Lab Sample ID: 240-57597-6 Date Sampled: 11/05/2015 0910
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111011.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.15 g |
| Analysis Date: | 11/11/2015 1319 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 83 |
| Di-n-octyl phthalate | | 9.4 | U | 9.4 | 83 |
| Fluoranthene | | 0.66 | U | 0.66 | 8.0 |
| Fluorene | | 0.63 | U | 0.63 | 8.0 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 8.0 |
| Hexachlorobutadiene | | 6.7 | U | 6.7 | 60 |
| Hexachlorocyclopentadiene | | 9.7 | U | 9.7 | 390 |
| Hexachloroethane | | 11 | U | 11 | 60 |
| Indeno[1,2,3-cd]pyrene | | 0.42 | U | 0.42 | 8.0 |
| Isophorone | | 15 | U | 15 | 60 |
| N-Nitrosodi-n-propylamine | | 7.5 | U | 7.5 | 60 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 60 |
| Naphthalene | | 0.98 | U | 0.98 | 8.0 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.87 | U | 0.87 | 8.0 |
| Phenol | | 8.7 | U | 8.7 | 60 |
| Pyrene | | 0.52 | U | 0.52 | 8.0 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 79 | | 36 - 110 |
| Phenol-d5 (Surr) | 60 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 62 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 62 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 68 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 33 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-7-14**

Lab Sample ID: 240-57597-7 Date Sampled: 11/05/2015 0912
Client Matrix: Solid % Moisture: 16.2 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111012.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.53 g |
| Analysis Date: | 11/11/2015 1344 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 61 |
| bis (2-chloroisopropyl) ether | | 12 | U | 12 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 400 |
| 2,4-Dinitrotoluene | | 21 | U | 21 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.55 | U | 0.55 | 61 |
| 2-Chlorophenol | | 9.9 | U | 9.9 | 61 |
| 2-Methylnaphthalene | | 0.61 | U | 0.61 | 8.1 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 10 | U | 10 | 61 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 21 | U | 21 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Nitroaniline | | 32 | U | 32 | 240 |
| 4-Nitrophenol | | 21 | U | 21 | 400 |
| Acenaphthene | | 0.92 | U | 0.92 | 8.1 |
| Acenaphthylene | | 0.42 | U | 0.42 | 8.1 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.95 | U | 0.95 | 8.1 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 15 | U | 15 | 120 |
| Benzo[a]anthracene | | 0.76 | U | 0.76 | 8.1 |
| Benzo[a]pyrene | | 0.78 | U | 0.78 | 8.1 |
| Benzo[b]fluoranthene | | 0.72 | U | 0.72 | 8.1 |
| Benzo[g,h,i]perylene | | 0.42 | U | 0.42 | 8.1 |
| Benzo[k]fluoranthene | | 0.82 | U | 0.82 | 8.1 |
| Bis(2-chloroethoxy)methane | | 27 | U | 27 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 32 | J | 23 | 85 |
| Butyl benzyl phthalate | | 12 | U | 12 | 85 |
| Caprolactam | | 45 | U | 45 | 400 |
| Carbazole | | 33 | U | 33 | 61 |
| Chrysene | | 1.3 | U | 1.3 | 8.1 |
| Dibenz(a,h)anthracene | | 0.80 | U | 0.80 | 8.1 |
| Dibenzofuran | | 0.80 | U | 0.80 | 61 |
| Diethyl phthalate | | 19 | U | 19 | 85 |
| Dimethyl phthalate | | 21 | U | 21 | 85 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-7-14**

Lab Sample ID: 240-57597-7 Date Sampled: 11/05/2015 0912
Client Matrix: Solid % Moisture: 16.2 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111012.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.53 g |
| Analysis Date: | 11/11/2015 1344 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 85 |
| Di-n-octyl phthalate | | 9.6 | U | 9.6 | 85 |
| Fluoranthene | | 0.67 | U | 0.67 | 8.1 |
| Fluorene | | 0.64 | U | 0.64 | 8.1 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 8.1 |
| Hexachlorobutadiene | | 6.8 | U | 6.8 | 61 |
| Hexachlorocyclopentadiene | | 9.8 | U | 9.8 | 400 |
| Hexachloroethane | | 11 | U | 11 | 61 |
| Indeno[1,2,3-cd]pyrene | | 0.42 | U | 0.42 | 8.1 |
| Isophorone | | 16 | U | 16 | 61 |
| N-Nitrosodi-n-propylamine | | 7.6 | U | 7.6 | 61 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 61 |
| Naphthalene | | 0.99 | U | 0.99 | 8.1 |
| Nitrobenzene | | 2.7 | U | 2.7 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.89 | U | 0.89 | 8.1 |
| Phenol | | 8.9 | U | 8.9 | 61 |
| Pyrene | | 0.53 | U | 0.53 | 8.1 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 490 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 80 | | 36 - 110 |
| Phenol-d5 (Surr) | 64 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 62 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 67 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 70 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 38 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-8-14**

Lab Sample ID: 240-57597-8 Date Sampled: 11/05/2015 0914
Client Matrix: Solid % Moisture: 18.0 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111013.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.20 g |
| Analysis Date: | 11/11/2015 1408 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 61 |
| bis (2-chloroisopropyl) ether | | 12 | U | 12 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 400 |
| 2,4-Dinitrotoluene | | 21 | U | 21 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.55 | U | 0.55 | 61 |
| 2-Chlorophenol | | 9.9 | U | 9.9 | 61 |
| 2-Methylnaphthalene | | 0.61 | U | 0.61 | 8.1 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 10 | U | 10 | 61 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 21 | U | 21 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 21 | U | 21 | 400 |
| Acenaphthene | | 0.92 | U | 0.92 | 8.1 |
| Acenaphthylene | | 0.42 | U | 0.42 | 8.1 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.94 | U | 0.94 | 8.1 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 15 | U | 15 | 120 |
| Benzo[a]anthracene | | 0.76 | U | 0.76 | 8.1 |
| Benzo[a]pyrene | | 0.78 | U | 0.78 | 8.1 |
| Benzo[b]fluoranthene | | 0.71 | U | 0.71 | 8.1 |
| Benzo[g,h,i]perylene | | 0.42 | U | 0.42 | 8.1 |
| Benzo[k]fluoranthene | | 0.82 | U | 0.82 | 8.1 |
| Bis(2-chloroethoxy)methane | | 27 | U | 27 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 23 | U | 23 | 85 |
| Butyl benzyl phthalate | | 12 | U | 12 | 85 |
| Caprolactam | | 45 | U | 45 | 400 |
| Carbazole | | 33 | U | 33 | 61 |
| Chrysene | | 1.3 | U | 1.3 | 8.1 |
| Dibenz(a,h)anthracene | | 0.80 | U | 0.80 | 8.1 |
| Dibenzofuran | | 0.80 | U | 0.80 | 61 |
| Diethyl phthalate | | 19 | U | 19 | 85 |
| Dimethyl phthalate | | 21 | U | 21 | 85 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-8-14**

Lab Sample ID: 240-57597-8 Date Sampled: 11/05/2015 0914
Client Matrix: Solid % Moisture: 18.0 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111013.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.20 g |
| Analysis Date: | 11/11/2015 1408 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 85 |
| Di-n-octyl phthalate | | 9.6 | U | 9.6 | 85 |
| Fluoranthene | | 0.67 | U | 0.67 | 8.1 |
| Fluorene | | 0.64 | U | 0.64 | 8.1 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 8.1 |
| Hexachlorobutadiene | | 6.8 | U | 6.8 | 61 |
| Hexachlorocyclopentadiene | | 9.8 | U | 9.8 | 400 |
| Hexachloroethane | | 11 | U | 11 | 61 |
| Indeno[1,2,3-cd]pyrene | | 0.42 | U | 0.42 | 8.1 |
| Isophorone | | 16 | U | 16 | 61 |
| N-Nitrosodi-n-propylamine | | 7.6 | U | 7.6 | 61 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 61 |
| Naphthalene | | 0.99 | U | 0.99 | 8.1 |
| Nitrobenzene | | 2.7 | U | 2.7 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.88 | U | 0.88 | 8.1 |
| Phenol | | 8.8 | U | 8.8 | 61 |
| Pyrene | | 0.53 | U | 0.53 | 8.1 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 70 | | 36 - 110 |
| Phenol-d5 (Surr) | 54 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 52 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 56 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 56 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 32 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-9-14**

Lab Sample ID: 240-57597-9 Date Sampled: 11/05/2015 0916
Client Matrix: Solid % Moisture: 19.1 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111014.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.67 g |
| Analysis Date: | 11/11/2015 1433 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.4 | U | 4.4 | 62 |
| bis (2-chloroisopropyl) ether | | 12 | U | 12 | 120 |
| 2,4,5-Trichlorophenol | | 31 | U | 31 | 190 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 190 |
| 2,4-Dichlorophenol | | 25 | U | 25 | 190 |
| 2,4-Dimethylphenol | | 25 | U | 25 | 190 |
| 2,4-Dinitrophenol | | 26 | U | 26 | 410 |
| 2,4-Dinitrotoluene | | 21 | U | 21 | 250 |
| 2,6-Dinitrotoluene | | 26 | U | 26 | 250 |
| 2-Chloronaphthalene | | 0.56 | U | 0.56 | 62 |
| 2-Chlorophenol | | 10 | U | 10 | 62 |
| 2-Methylnaphthalene | | 0.62 | U | 0.62 | 8.3 |
| 2-Methylphenol | | 14 | U | 14 | 250 |
| 2-Nitroaniline | | 11 | U | 11 | 250 |
| 2-Nitrophenol | | 10 | U | 10 | 62 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 20 | U | 20 | 250 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 190 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 62 |
| 4-Chloro-3-methylphenol | | 26 | U | 26 | 190 |
| 4-Chloroaniline | | 21 | U | 21 | 190 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 62 |
| 4-Nitroaniline | | 32 | U | 32 | 250 |
| 4-Nitrophenol | | 21 | U | 21 | 410 |
| Acenaphthene | | 0.95 | U | 0.95 | 8.3 |
| Acenaphthylene | | 0.44 | U | 0.44 | 8.3 |
| Acetophenone | | 13 | J | 11 | 120 |
| Anthracene | | 0.97 | U | 0.97 | 8.3 |
| Atrazine | | 11 | U | 11 | 250 |
| Benzaldehyde | | 15 | U | 15 | 120 |
| Benzo[a]anthracene | | 0.79 | U | 0.79 | 8.3 |
| Benzo[a]pyrene | | 0.80 | U | 0.80 | 8.3 |
| Benzo[b]fluoranthene | | 0.74 | U | 0.74 | 8.3 |
| Benzo[g,h,i]perylene | | 0.44 | U | 0.44 | 8.3 |
| Benzo[k]fluoranthene | | 0.85 | U | 0.85 | 8.3 |
| Bis(2-chloroethoxy)methane | | 27 | U | 27 | 120 |
| Bis(2-chloroethyl)ether | | 2.5 | U | 2.5 | 120 |
| Bis(2-ethylhexyl) phthalate | | 40 | J | 24 | 87 |
| Butyl benzyl phthalate | | 12 | U | 12 | 87 |
| Caprolactam | | 46 | U | 46 | 410 |
| Carbazole | | 34 | U | 34 | 62 |
| Chrysene | | 1.4 | U | 1.4 | 8.3 |
| Dibenz(a,h)anthracene | | 0.82 | U | 0.82 | 8.3 |
| Dibenzofuran | | 0.82 | U | 0.82 | 62 |
| Diethyl phthalate | | 20 | U | 20 | 87 |
| Dimethyl phthalate | | 21 | U | 21 | 87 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-9-14**

Lab Sample ID: 240-57597-9 Date Sampled: 11/05/2015 0916
Client Matrix: Solid % Moisture: 19.1 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111014.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.67 g |
| Analysis Date: | 11/11/2015 1433 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 19 | U | 19 | 87 |
| Di-n-octyl phthalate | | 9.9 | U | 9.9 | 87 |
| Fluoranthene | | 0.69 | U | 0.69 | 8.3 |
| Fluorene | | 0.66 | U | 0.66 | 8.3 |
| Hexachlorobenzene | | 2.6 | U | 2.6 | 8.3 |
| Hexachlorobutadiene | | 7.0 | U | 7.0 | 62 |
| Hexachlorocyclopentadiene | | 10 | U | 10 | 410 |
| Hexachloroethane | | 11 | U | 11 | 62 |
| Indeno[1,2,3-cd]pyrene | | 0.44 | U | 0.44 | 8.3 |
| Isophorone | | 16 | U | 16 | 62 |
| N-Nitrosodi-n-propylamine | | 7.9 | U | 7.9 | 62 |
| N-Nitrosodiphenylamine | | 26 | U | 26 | 62 |
| Naphthalene | | 1.0 | U | 1.0 | 8.3 |
| Nitrobenzene | | 2.7 | U | 2.7 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 190 |
| Phenanthrene | | 0.91 | U | 0.91 | 8.3 |
| Phenol | | 9.1 | U | 9.1 | 62 |
| Pyrene | | 0.55 | U | 0.55 | 8.3 |
| 3 & 4 Methylphenol | | 25 | U | 25 | 500 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 87 | | 36 - 110 |
| Phenol-d5 (Surr) | 68 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 68 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 71 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 73 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 45 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-10-14**

Lab Sample ID: 240-57597-10 Date Sampled: 11/05/2015 0918
Client Matrix: Solid % Moisture: 13.4 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111015.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.78 g |
| Analysis Date: | 11/11/2015 1457 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 58 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 29 | U | 29 | 170 |
| 2,4,6-Trichlorophenol | | 10 | U | 10 | 170 |
| 2,4-Dichlorophenol | | 23 | U | 23 | 170 |
| 2,4-Dimethylphenol | | 23 | U | 23 | 170 |
| 2,4-Dinitrophenol | | 24 | U | 24 | 380 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 230 |
| 2,6-Dinitrotoluene | | 24 | U | 24 | 230 |
| 2-Chloronaphthalene | | 0.52 | U | 0.52 | 58 |
| 2-Chlorophenol | | 9.5 | U | 9.5 | 58 |
| 2-Methylnaphthalene | | 0.58 | U | 0.58 | 7.8 |
| 2-Methylphenol | | 13 | U | 13 | 230 |
| 2-Nitroaniline | | 11 | U | 11 | 230 |
| 2-Nitrophenol | | 9.7 | U | 9.7 | 58 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 230 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 170 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Chloro-3-methylphenol | | 24 | U | 24 | 170 |
| 4-Chloroaniline | | 20 | U | 20 | 170 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Nitroaniline | | 30 | U | 30 | 230 |
| 4-Nitrophenol | | 20 | U | 20 | 380 |
| Acenaphthene | | 0.88 | U | 0.88 | 7.8 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.8 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.91 | U | 0.91 | 7.8 |
| Atrazine | | 11 | U | 11 | 230 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.73 | U | 0.73 | 7.8 |
| Benzo[a]pyrene | | 0.74 | U | 0.74 | 7.8 |
| Benzo[b]fluoranthene | | 0.69 | U | 0.69 | 7.8 |
| Benzo[g,h,i]perylene | | 0.41 | U | 0.41 | 7.8 |
| Benzo[k]fluoranthene | | 0.79 | U | 0.79 | 7.8 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.3 | U | 2.3 | 120 |
| Bis(2-ethylhexyl) phthalate | | 29 | J | 22 | 81 |
| Butyl benzyl phthalate | | 12 | U | 12 | 81 |
| Caprolactam | | 43 | U | 43 | 380 |
| Carbazole | | 31 | U | 31 | 58 |
| Chrysene | | 1.3 | U | 1.3 | 7.8 |
| Dibenz(a,h)anthracene | | 0.77 | U | 0.77 | 7.8 |
| Dibenzofuran | | 0.77 | U | 0.77 | 58 |
| Diethyl phthalate | | 19 | U | 19 | 81 |
| Dimethyl phthalate | | 20 | U | 20 | 81 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-10-14**

Lab Sample ID: 240-57597-10 Date Sampled: 11/05/2015 0918
Client Matrix: Solid % Moisture: 13.4 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-205786 | Lab File ID: | 51111015.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.78 g |
| Analysis Date: | 11/11/2015 1457 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 17 | U | 17 | 81 |
| Di-n-octyl phthalate | | 9.2 | U | 9.2 | 81 |
| Fluoranthene | | 0.64 | U | 0.64 | 7.8 |
| Fluorene | | 0.62 | U | 0.62 | 7.8 |
| Hexachlorobenzene | | 2.4 | U | 2.4 | 7.8 |
| Hexachlorobutadiene | | 6.5 | U | 6.5 | 58 |
| Hexachlorocyclopentadiene | | 9.4 | U | 9.4 | 380 |
| Hexachloroethane | | 10 | U | 10 | 58 |
| Indeno[1,2,3-cd]pyrene | | 0.41 | U | 0.41 | 7.8 |
| Isophorone | | 15 | U | 15 | 58 |
| N-Nitrosodi-n-propylamine | | 7.3 | U | 7.3 | 58 |
| N-Nitrosodiphenylamine | | 24 | U | 24 | 58 |
| Naphthalene | | 0.95 | U | 0.95 | 7.8 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 170 |
| Phenanthrene | | 0.85 | U | 0.85 | 7.8 |
| Phenol | | 8.5 | U | 8.5 | 58 |
| Pyrene | | 0.51 | U | 0.51 | 7.8 |
| 3 & 4 Methylphenol | | 23 | U | 23 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 78 | | 36 - 110 |
| Phenol-d5 (Surr) | 61 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 60 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 59 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 64 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 40 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-11-14**

Lab Sample ID: 240-57597-11 Date Sampled: 11/05/2015 0920
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111019.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.86 g |
| Analysis Date: | 11/11/2015 1450 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 60 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 400 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.54 | U | 0.54 | 60 |
| 2-Chlorophenol | | 9.9 | U | 9.9 | 60 |
| 2-Methylnaphthalene | | 0.60 | U | 0.60 | 8.0 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 10 | U | 10 | 60 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 60 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 60 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 400 |
| Acenaphthene | | 0.91 | U | 0.91 | 8.0 |
| Acenaphthylene | | 0.42 | U | 0.42 | 8.0 |
| Acetophenone | | 38 | J | 11 | 120 |
| Anthracene | | 0.94 | U | 0.94 | 8.0 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 19 | J | 14 | 120 |
| Benzo[a]anthracene | | 0.76 | U | 0.76 | 8.0 |
| Benzo[a]pyrene | | 0.77 | U | 0.77 | 8.0 |
| Benzo[b]fluoranthene | | 0.71 | U | 0.71 | 8.0 |
| Benzo[g,h,i]perylene | | 0.42 | U | 0.42 | 8.0 |
| Benzo[k]fluoranthene | | 0.82 | U | 0.82 | 8.0 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 23 | U | 23 | 84 |
| Butyl benzyl phthalate | | 12 | U | 12 | 84 |
| Caprolactam | | 45 | U | 45 | 400 |
| Carbazole | | 32 | U | 32 | 60 |
| Chrysene | | 1.3 | U | 1.3 | 8.0 |
| Dibenz(a,h)anthracene | | 0.79 | U | 0.79 | 8.0 |
| Dibenzofuran | | 0.79 | U | 0.79 | 60 |
| Diethyl phthalate | | 19 | U | 19 | 84 |
| Dimethyl phthalate | | 20 | U | 20 | 84 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-11-14**

Lab Sample ID: 240-57597-11 Date Sampled: 11/05/2015 0920
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111019.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.86 g |
| Analysis Date: | 11/11/2015 1450 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 84 |
| Di-n-octyl phthalate | | 9.5 | U | 9.5 | 84 |
| Fluoranthene | | 0.66 | U | 0.66 | 8.0 |
| Fluorene | | 0.64 | U | 0.64 | 8.0 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 8.0 |
| Hexachlorobutadiene | | 6.7 | U | 6.7 | 60 |
| Hexachlorocyclopentadiene | | 9.7 | U | 9.7 | 400 |
| Hexachloroethane | | 11 | U | 11 | 60 |
| Indeno[1,2,3-cd]pyrene | | 0.42 | U | 0.42 | 8.0 |
| Isophorone | | 16 | U | 16 | 60 |
| N-Nitrosodi-n-propylamine | | 7.6 | U | 7.6 | 60 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 60 |
| Naphthalene | | 0.99 | U | 0.99 | 8.0 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.88 | U | 0.88 | 8.0 |
| Phenol | | 8.8 | U | 8.8 | 60 |
| Pyrene | | 0.53 | U | 0.53 | 8.0 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 88 | | 36 - 110 |
| Phenol-d5 (Surr) | 80 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 80 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 78 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 73 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 40 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-12-14**

Lab Sample ID: 240-57597-12 Date Sampled: 11/05/2015 0922
Client Matrix: Solid % Moisture: 15.9 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111020.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.84 g |
| Analysis Date: | 11/11/2015 1514 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 60 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.54 | U | 0.54 | 60 |
| 2-Chlorophenol | | 9.8 | U | 9.8 | 60 |
| 2-Methylnaphthalene | | 0.60 | U | 0.60 | 8.0 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.9 | U | 9.9 | 60 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 60 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 60 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.91 | U | 0.91 | 8.0 |
| Acenaphthylene | | 0.42 | U | 0.42 | 8.0 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.93 | U | 0.93 | 8.0 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.75 | U | 0.75 | 8.0 |
| Benzo[a]pyrene | | 0.77 | U | 0.77 | 8.0 |
| Benzo[b]fluoranthene | | 0.71 | U | 0.71 | 8.0 |
| Benzo[g,h,i]perylene | | 0.42 | U | 0.42 | 8.0 |
| Benzo[k]fluoranthene | | 0.81 | U | 0.81 | 8.0 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 23 | U | 23 | 84 |
| Butyl benzyl phthalate | | 12 | U | 12 | 84 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 60 |
| Chrysene | | 1.3 | U | 1.3 | 8.0 |
| Dibenz(a,h)anthracene | | 0.79 | U | 0.79 | 8.0 |
| Dibenzofuran | | 0.79 | U | 0.79 | 60 |
| Diethyl phthalate | | 19 | U | 19 | 84 |
| Dimethyl phthalate | | 20 | U | 20 | 84 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-12-14**

Lab Sample ID: 240-57597-12 Date Sampled: 11/05/2015 0922
Client Matrix: Solid % Moisture: 15.9 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111020.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.84 g |
| Analysis Date: | 11/11/2015 1514 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 84 |
| Di-n-octyl phthalate | | 9.4 | U | 9.4 | 84 |
| Fluoranthene | | 0.66 | U | 0.66 | 8.0 |
| Fluorene | | 0.63 | U | 0.63 | 8.0 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 8.0 |
| Hexachlorobutadiene | | 6.7 | U | 6.7 | 60 |
| Hexachlorocyclopentadiene | | 9.7 | U | 9.7 | 390 |
| Hexachloroethane | | 11 | U | 11 | 60 |
| Indeno[1,2,3-cd]pyrene | | 0.42 | U | 0.42 | 8.0 |
| Isophorone | | 16 | U | 16 | 60 |
| N-Nitrosodi-n-propylamine | | 7.5 | U | 7.5 | 60 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 60 |
| Naphthalene | | 0.98 | U | 0.98 | 8.0 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.87 | U | 0.87 | 8.0 |
| Phenol | | 8.7 | U | 8.7 | 60 |
| Pyrene | | 0.53 | U | 0.53 | 8.0 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 72 | | 36 - 110 |
| Phenol-d5 (Surr) | 64 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 67 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 67 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 63 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 32 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-13-14**

Lab Sample ID: 240-57597-13 Date Sampled: 11/05/2015 0924
Client Matrix: Solid % Moisture: 14.7 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111021.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.37 g |
| Analysis Date: | 11/11/2015 1539 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 58 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 29 | U | 29 | 170 |
| 2,4,6-Trichlorophenol | | 10 | U | 10 | 170 |
| 2,4-Dichlorophenol | | 23 | U | 23 | 170 |
| 2,4-Dimethylphenol | | 23 | U | 23 | 170 |
| 2,4-Dinitrophenol | | 24 | U | 24 | 380 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 230 |
| 2,6-Dinitrotoluene | | 24 | U | 24 | 230 |
| 2-Chloronaphthalene | | 0.52 | U | 0.52 | 58 |
| 2-Chlorophenol | | 9.5 | U | 9.5 | 58 |
| 2-Methylnaphthalene | | 0.58 | U | 0.58 | 7.7 |
| 2-Methylphenol | | 13 | U | 13 | 230 |
| 2-Nitroaniline | | 11 | U | 11 | 230 |
| 2-Nitrophenol | | 9.6 | U | 9.6 | 58 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 230 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 170 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Chloro-3-methylphenol | | 24 | U | 24 | 170 |
| 4-Chloroaniline | | 20 | U | 20 | 170 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Nitroaniline | | 30 | U | 30 | 230 |
| 4-Nitrophenol | | 20 | U | 20 | 380 |
| Acenaphthene | | 0.88 | U | 0.88 | 7.7 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.7 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.90 | U | 0.90 | 7.7 |
| Atrazine | | 11 | U | 11 | 230 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.73 | U | 0.73 | 7.7 |
| Benzo[a]pyrene | | 0.74 | U | 0.74 | 7.7 |
| Benzo[b]fluoranthene | | 0.68 | U | 0.68 | 7.7 |
| Benzo[g,h,i]perylene | | 0.41 | U | 0.41 | 7.7 |
| Benzo[k]fluoranthene | | 0.79 | U | 0.79 | 7.7 |
| Bis(2-chloroethoxy)methane | | 25 | U | 25 | 120 |
| Bis(2-chloroethyl)ether | | 2.3 | U | 2.3 | 120 |
| Bis(2-ethylhexyl) phthalate | | 60 | J | 22 | 81 |
| Butyl benzyl phthalate | | 12 | U | 12 | 81 |
| Caprolactam | | 43 | U | 43 | 380 |
| Carbazole | | 31 | U | 31 | 58 |
| Chrysene | | 1.3 | U | 1.3 | 7.7 |
| Dibenz(a,h)anthracene | | 0.76 | U | 0.76 | 7.7 |
| Dibenzofuran | | 0.76 | U | 0.76 | 58 |
| Diethyl phthalate | | 19 | U | 19 | 81 |
| Dimethyl phthalate | | 20 | U | 20 | 81 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-13-14**

Lab Sample ID: 240-57597-13 Date Sampled: 11/05/2015 0924
Client Matrix: Solid % Moisture: 14.7 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111021.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.37 g |
| Analysis Date: | 11/11/2015 1539 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | J | 17 | 81 |
| Di-n-octyl phthalate | | 9.1 | U | 9.1 | 81 |
| Fluoranthene | | 0.64 | U | 0.64 | 7.7 |
| Fluorene | | 0.61 | U | 0.61 | 7.7 |
| Hexachlorobenzene | | 2.4 | U | 2.4 | 7.7 |
| Hexachlorobutadiene | | 6.5 | U | 6.5 | 58 |
| Hexachlorocyclopentadiene | | 9.4 | U | 9.4 | 380 |
| Hexachloroethane | | 10 | U | 10 | 58 |
| Indeno[1,2,3-cd]pyrene | | 0.41 | U | 0.41 | 7.7 |
| Isophorone | | 15 | U | 15 | 58 |
| N-Nitrosodi-n-propylamine | | 7.3 | U | 7.3 | 58 |
| N-Nitrosodiphenylamine | | 24 | U | 24 | 58 |
| Naphthalene | | 0.95 | U | 0.95 | 7.7 |
| Nitrobenzene | | 2.5 | U | 2.5 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 170 |
| Phenanthrene | | 0.85 | U | 0.85 | 7.7 |
| Phenol | | 8.5 | U | 8.5 | 58 |
| Pyrene | | 0.51 | U | 0.51 | 7.7 |
| 3 & 4 Methylphenol | | 23 | U | 23 | 460 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 77 | | 36 - 110 |
| Phenol-d5 (Surr) | 57 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 65 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 65 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 59 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 33 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-14-14**

Lab Sample ID: 240-57597-14 Date Sampled: 11/05/2015 0926
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111022.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.64 g |
| Analysis Date: | 11/11/2015 1603 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 59 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.53 | U | 0.53 | 59 |
| 2-Chlorophenol | | 9.7 | U | 9.7 | 59 |
| 2-Methylnaphthalene | | 0.59 | U | 0.59 | 7.9 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.8 | U | 9.8 | 59 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.90 | U | 0.90 | 7.9 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.9 |
| Acetophenone | | 11 | J | 11 | 120 |
| Anthracene | | 0.92 | U | 0.92 | 7.9 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.75 | U | 0.75 | 7.9 |
| Benzo[a]pyrene | | 0.76 | U | 0.76 | 7.9 |
| Benzo[b]fluoranthene | | 0.70 | U | 0.70 | 7.9 |
| Benzo[g,h,i]perylene | | 0.41 | U | 0.41 | 7.9 |
| Benzo[k]fluoranthene | | 0.80 | U | 0.80 | 7.9 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 22 | U | 22 | 83 |
| Butyl benzyl phthalate | | 12 | U | 12 | 83 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 59 |
| Chrysene | | 1.3 | U | 1.3 | 7.9 |
| Dibenz(a,h)anthracene | | 0.78 | U | 0.78 | 7.9 |
| Dibenzofuran | | 0.78 | U | 0.78 | 59 |
| Diethyl phthalate | | 19 | U | 19 | 83 |
| Dimethyl phthalate | | 20 | U | 20 | 83 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-14-14**

Lab Sample ID: 240-57597-14 Date Sampled: 11/05/2015 0926
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111022.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.64 g |
| Analysis Date: | 11/11/2015 1603 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 83 |
| Di-n-octyl phthalate | | 9.4 | U | 9.4 | 83 |
| Fluoranthene | | 0.65 | U | 0.65 | 7.9 |
| Fluorene | | 0.63 | U | 0.63 | 7.9 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.9 |
| Hexachlorobutadiene | | 6.6 | U | 6.6 | 59 |
| Hexachlorocyclopentadiene | | 9.6 | U | 9.6 | 390 |
| Hexachloroethane | | 11 | U | 11 | 59 |
| Indeno[1,2,3-cd]pyrene | | 0.41 | U | 0.41 | 7.9 |
| Isophorone | | 15 | U | 15 | 59 |
| N-Nitrosodi-n-propylamine | | 7.5 | U | 7.5 | 59 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 59 |
| Naphthalene | | 0.97 | U | 0.97 | 7.9 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.86 | U | 0.86 | 7.9 |
| Phenol | | 8.6 | U | 8.6 | 59 |
| Pyrene | | 0.52 | U | 0.52 | 7.9 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 79 | | 36 - 110 |
| Phenol-d5 (Surr) | 62 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 73 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 70 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 68 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 28 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-DUP**

Lab Sample ID: 240-57597-15 Date Sampled: 11/05/2015 0000
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111023.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.09 g |
| Analysis Date: | 11/11/2015 1628 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 58 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 29 | U | 29 | 170 |
| 2,4,6-Trichlorophenol | | 10 | U | 10 | 170 |
| 2,4-Dichlorophenol | | 23 | U | 23 | 170 |
| 2,4-Dimethylphenol | | 23 | U | 23 | 170 |
| 2,4-Dinitrophenol | | 24 | U | 24 | 380 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 230 |
| 2,6-Dinitrotoluene | | 24 | U | 24 | 230 |
| 2-Chloronaphthalene | | 0.52 | U | 0.52 | 58 |
| 2-Chlorophenol | | 9.6 | U | 9.6 | 58 |
| 2-Methylnaphthalene | | 0.58 | U | 0.58 | 7.8 |
| 2-Methylphenol | | 13 | U | 13 | 230 |
| 2-Nitroaniline | | 11 | U | 11 | 230 |
| 2-Nitrophenol | | 9.7 | U | 9.7 | 58 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 230 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 170 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Chloro-3-methylphenol | | 24 | U | 24 | 170 |
| 4-Chloroaniline | | 20 | U | 20 | 170 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Nitroaniline | | 30 | U | 30 | 230 |
| 4-Nitrophenol | | 20 | U | 20 | 380 |
| Acenaphthene | | 0.89 | U | 0.89 | 7.8 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.8 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.91 | U | 0.91 | 7.8 |
| Atrazine | | 11 | U | 11 | 230 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.73 | U | 0.73 | 7.8 |
| Benzo[a]pyrene | | 0.75 | U | 0.75 | 7.8 |
| Benzo[b]fluoranthene | | 0.69 | U | 0.69 | 7.8 |
| Benzo[g,h,i]perylene | | 0.41 | U | 0.41 | 7.8 |
| Benzo[k]fluoranthene | | 0.79 | U | 0.79 | 7.8 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.3 | U | 2.3 | 120 |
| Bis(2-ethylhexyl) phthalate | | 22 | J | 22 | 82 |
| Butyl benzyl phthalate | | 12 | U | 12 | 82 |
| Caprolactam | | 43 | U | 43 | 380 |
| Carbazole | | 31 | U | 31 | 58 |
| Chrysene | | 1.3 | U | 1.3 | 7.8 |
| Dibenz(a,h)anthracene | | 0.77 | U | 0.77 | 7.8 |
| Dibenzofuran | | 0.77 | U | 0.77 | 58 |
| Diethyl phthalate | | 19 | U | 19 | 82 |
| Dimethyl phthalate | | 20 | U | 20 | 82 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-DUP**

Lab Sample ID: 240-57597-15 Date Sampled: 11/05/2015 0000
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-205866 | Lab File ID: | 51111023.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.09 g |
| Analysis Date: | 11/11/2015 1628 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 17 | U | 17 | 82 |
| Di-n-octyl phthalate | | 9.2 | U | 9.2 | 82 |
| Fluoranthene | | 0.64 | U | 0.64 | 7.8 |
| Fluorene | | 0.62 | U | 0.62 | 7.8 |
| Hexachlorobenzene | | 2.4 | U | 2.4 | 7.8 |
| Hexachlorobutadiene | | 6.5 | U | 6.5 | 58 |
| Hexachlorocyclopentadiene | | 9.4 | U | 9.4 | 380 |
| Hexachloroethane | | 10 | U | 10 | 58 |
| Indeno[1,2,3-cd]pyrene | | 0.41 | U | 0.41 | 7.8 |
| Isophorone | | 15 | U | 15 | 58 |
| N-Nitrosodi-n-propylamine | | 7.3 | U | 7.3 | 58 |
| N-Nitrosodiphenylamine | | 24 | U | 24 | 58 |
| Naphthalene | | 0.96 | U | 0.96 | 7.8 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 170 |
| Phenanthrene | | 0.85 | U | 0.85 | 7.8 |
| Phenol | | 8.5 | U | 8.5 | 58 |
| Pyrene | | 0.51 | U | 0.51 | 7.8 |
| 3 & 4 Methylphenol | | 23 | U | 23 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 81 | | 36 - 110 |
| Phenol-d5 (Surr) | 60 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 74 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 71 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 68 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 24 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-1-14**

Lab Sample ID: 240-57597-1 Date Sampled: 11/05/2015 0900
Client Matrix: Solid % Moisture: 15.7 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 4.98 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1108 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 78 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-2-14**

Lab Sample ID: 240-57597-2 Date Sampled: 11/05/2015 0902
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 4.98 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1145 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 54 | U | 54 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 79 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-3-14**

Lab Sample ID: 240-57597-3 Date Sampled: 11/05/2015 0904
Client Matrix: Solid % Moisture: 17.8 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 4.99 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1223 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 56 | U | 56 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 77 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-4-14**

Lab Sample ID: 240-57597-4 Date Sampled: 11/05/2015 0906
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 4.99 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1300 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 71 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-5-14**

Lab Sample ID: 240-57597-5 Date Sampled: 11/05/2015 0908
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206404 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206444 | Initial Weight/Volume: | 4.98 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 0917 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/12/2015 0559 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 77 | J | 54 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 71 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-6-14**

Lab Sample ID: 240-57597-6 Date Sampled: 11/05/2015 0910
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 5.01 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1413 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 76 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-7-14**

Lab Sample ID: 240-57597-7 Date Sampled: 11/05/2015 0912
Client Matrix: Solid % Moisture: 16.2 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 4.97 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1450 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 78 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-8-14**

Lab Sample ID: 240-57597-8 Date Sampled: 11/05/2015 0914
Client Matrix: Solid % Moisture: 18.0 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 4.98 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1527 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 56 | U | 56 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 73 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-9-14**

Lab Sample ID: 240-57597-9 Date Sampled: 11/05/2015 0916
Client Matrix: Solid % Moisture: 19.1 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 5.00 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1719 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 57 | U | 57 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 78 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-10-14**

Lab Sample ID: 240-57597-10 Date Sampled: 11/05/2015 0918
Client Matrix: Solid % Moisture: 13.4 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 5.02 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1757 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 53 | U | 53 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 79 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-11-14**

Lab Sample ID: 240-57597-11 Date Sampled: 11/05/2015 0920
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 5.03 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1834 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 79 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-12-14**

Lab Sample ID: 240-57597-12 Date Sampled: 11/05/2015 0922
Client Matrix: Solid % Moisture: 15.9 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 5.04 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1912 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 54 | U | 54 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 78 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID: F-VB-UST-13-14**

Lab Sample ID: 240-57597-13 Date Sampled: 11/05/2015 0924
Client Matrix: Solid % Moisture: 14.7 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 4.99 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1948 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 54 | U | 54 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 75 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-14-14**

Lab Sample ID: 240-57597-14 Date Sampled: 11/05/2015 0926
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 5.02 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 2025 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 54 | U | 54 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 80 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-DUP**

Lab Sample ID: 240-57597-15 Date Sampled: 11/05/2015 0000
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-206227 | Initial Weight/Volume: | 5.00 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 2103 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 54 | U F2 | 54 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 74 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-1-14**

Lab Sample ID: 240-57597-1 Date Sampled: 11/05/2015 0900
Client Matrix: Solid % Moisture: 15.7 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 30.49 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/10/2015 2007 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 16 | J B | 7.5 | 19 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 62 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-2-14**

Lab Sample ID: 240-57597-2 Date Sampled: 11/05/2015 0902
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 29.69 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/10/2015 2038 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 10 | J B | 7.6 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 62 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-3-14**

Lab Sample ID: 240-57597-3 Date Sampled: 11/05/2015 0904
Client Matrix: Solid % Moisture: 17.8 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 30.32 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/10/2015 2110 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.7 | U | 7.7 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 59 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-4-14**

Lab Sample ID: 240-57597-4 Date Sampled: 11/05/2015 0906
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 30.31 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/10/2015 2141 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.6 | J B | 7.5 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 60 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-5-14**

Lab Sample ID: 240-57597-5 Date Sampled: 11/05/2015 0908
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 30.01 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/10/2015 2243 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.5 | U | 7.5 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 56 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-6-14**

Lab Sample ID: 240-57597-6 Date Sampled: 11/05/2015 0910
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 29.79 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/10/2015 2314 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.7 | U | 7.7 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 60 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-7-14**

Lab Sample ID: 240-57597-7 Date Sampled: 11/05/2015 0912
Client Matrix: Solid % Moisture: 16.2 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 29.99 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/10/2015 2346 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 16 | J B | 7.6 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 65 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-8-14**

Lab Sample ID: 240-57597-8 Date Sampled: 11/05/2015 0914
Client Matrix: Solid % Moisture: 18.0 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 30.45 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 0017 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.7 | U | 7.7 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 65 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-9-14**

Lab Sample ID: 240-57597-9 Date Sampled: 11/05/2015 0916
Client Matrix: Solid % Moisture: 19.1 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 30.38 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 0048 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.8 | U | 7.8 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 57 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-10-14**

Lab Sample ID: 240-57597-10 Date Sampled: 11/05/2015 0918
Client Matrix: Solid % Moisture: 13.4 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 30.04 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 0119 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 13 | J B | 7.4 | 19 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 64 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-11-14**

Lab Sample ID: 240-57597-11 Date Sampled: 11/05/2015 0920
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 29.60 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 0150 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 9.0 | J B | 7.8 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 64 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-12-14**

Lab Sample ID: 240-57597-12 Date Sampled: 11/05/2015 0922
Client Matrix: Solid % Moisture: 15.9 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206356 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 29.52 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1903 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 30 | B | 7.7 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 73 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-13-14**

Lab Sample ID: 240-57597-13 Date Sampled: 11/05/2015 0924
Client Matrix: Solid % Moisture: 14.7 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206356 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 29.73 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 1934 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 13 | J B | 7.6 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 68 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-14-14**

Lab Sample ID: 240-57597-14 Date Sampled: 11/05/2015 0926
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206356 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 30.43 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 2006 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 11 | J B | 7.4 | 19 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 61 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-DUP**

Lab Sample ID: 240-57597-15 Date Sampled: 11/05/2015 0000
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206356 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-205788 | Initial Weight/Volume: | 29.98 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/11/2015 2037 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0940 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 8.2 | J B | 7.5 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 65 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-1-14**

Lab Sample ID: 240-57597-1 Date Sampled: 11/05/2015 0900
Client Matrix: Solid % Moisture: 15.7 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 29.54 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1328 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 40 |
| Aroclor-1221 | | 19 | U | 19 | 40 |
| Aroclor-1232 | | 24 | U | 24 | 40 |
| Aroclor-1242 | | 13 | U | 13 | 40 |
| Aroclor-1248 | | 9.6 | U | 9.6 | 40 |
| Aroclor-1254 | | 17 | U | 17 | 40 |
| Aroclor-1260 | | 11 | U | 11 | 40 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 78 | | 13 - 134 |
| DCB Decachlorobiphenyl | 87 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-2-14**

Lab Sample ID: 240-57597-2 Date Sampled: 11/05/2015 0902
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.07 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1345 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 23 | U | 23 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.4 | U | 9.4 | 39 |
| Aroclor-1254 | | 16 | U | 16 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 84 | | 13 - 134 |
| DCB Decachlorobiphenyl | 100 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-3-14**

Lab Sample ID: 240-57597-3 Date Sampled: 11/05/2015 0904
Client Matrix: Solid % Moisture: 17.8 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.14 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1401 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 15 | U | 15 | 40 |
| Aroclor-1221 | | 19 | U | 19 | 40 |
| Aroclor-1232 | | 24 | U | 24 | 40 |
| Aroclor-1242 | | 13 | U | 13 | 40 |
| Aroclor-1248 | | 9.7 | U | 9.7 | 40 |
| Aroclor-1254 | | 17 | U | 17 | 40 |
| Aroclor-1260 | | 11 | U | 11 | 40 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 68 | | 13 - 134 |
| DCB Decachlorobiphenyl | 95 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-4-14**

Lab Sample ID: 240-57597-4 Date Sampled: 11/05/2015 0906
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 29.54 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1417 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 40 |
| Aroclor-1221 | | 19 | U | 19 | 40 |
| Aroclor-1232 | | 24 | U | 24 | 40 |
| Aroclor-1242 | | 13 | U | 13 | 40 |
| Aroclor-1248 | | 9.6 | U | 9.6 | 40 |
| Aroclor-1254 | | 17 | U | 17 | 40 |
| Aroclor-1260 | | 11 | U | 11 | 40 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 75 | | 13 - 134 |
| DCB Decachlorobiphenyl | 104 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-5-14**

Lab Sample ID: 240-57597-5 Date Sampled: 11/05/2015 0908
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.48 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1433 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 38 |
| Aroclor-1221 | | 19 | U | 19 | 38 |
| Aroclor-1232 | | 23 | U | 23 | 38 |
| Aroclor-1242 | | 13 | U | 13 | 38 |
| Aroclor-1248 | | 9.3 | U | 9.3 | 38 |
| Aroclor-1254 | | 16 | U | 16 | 38 |
| Aroclor-1260 | | 10 | U | 10 | 38 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 73 | | 13 - 134 |
| DCB Decachlorobiphenyl | 91 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-6-14**

Lab Sample ID: 240-57597-6 Date Sampled: 11/05/2015 0910
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 29.87 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1450 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 40 |
| Aroclor-1221 | | 19 | U | 19 | 40 |
| Aroclor-1232 | | 24 | U | 24 | 40 |
| Aroclor-1242 | | 13 | U | 13 | 40 |
| Aroclor-1248 | | 9.6 | U | 9.6 | 40 |
| Aroclor-1254 | | 17 | U | 17 | 40 |
| Aroclor-1260 | | 11 | U | 11 | 40 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 76 | | 13 - 134 |
| DCB Decachlorobiphenyl | 109 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-7-14**

Lab Sample ID: 240-57597-7 Date Sampled: 11/05/2015 0912
Client Matrix: Solid % Moisture: 16.2 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.04 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1506 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 24 | U | 24 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.5 | U | 9.5 | 39 |
| Aroclor-1254 | | 17 | U | 17 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 76 | | 13 - 134 |
| DCB Decachlorobiphenyl | 98 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-8-14**

Lab Sample ID: 240-57597-8 Date Sampled: 11/05/2015 0914
Client Matrix: Solid % Moisture: 18.0 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 29.74 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1523 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 15 | U | 15 | 41 |
| Aroclor-1221 | | 20 | U | 20 | 41 |
| Aroclor-1232 | | 25 | U | 25 | 41 |
| Aroclor-1242 | | 14 | U | 14 | 41 |
| Aroclor-1248 | | 9.8 | U | 9.8 | 41 |
| Aroclor-1254 | | 17 | U | 17 | 41 |
| Aroclor-1260 | | 11 | U | 11 | 41 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 72 | | 13 - 134 |
| DCB Decachlorobiphenyl | 90 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-9-14**

Lab Sample ID: 240-57597-9 Date Sampled: 11/05/2015 0916
Client Matrix: Solid % Moisture: 19.1 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.07 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1539 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 15 | U | 15 | 41 |
| Aroclor-1221 | | 20 | U | 20 | 41 |
| Aroclor-1232 | | 25 | U | 25 | 41 |
| Aroclor-1242 | | 14 | U | 14 | 41 |
| Aroclor-1248 | | 9.9 | U | 9.9 | 41 |
| Aroclor-1254 | | 17 | U | 17 | 41 |
| Aroclor-1260 | | 11 | U | 11 | 41 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 77 | | 13 - 134 |
| DCB Decachlorobiphenyl | 95 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-10-14**

Lab Sample ID: 240-57597-10 Date Sampled: 11/05/2015 0918
Client Matrix: Solid % Moisture: 13.4 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 29.97 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1555 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 38 |
| Aroclor-1221 | | 19 | U | 19 | 38 |
| Aroclor-1232 | | 23 | U | 23 | 38 |
| Aroclor-1242 | | 13 | U | 13 | 38 |
| Aroclor-1248 | | 9.3 | U | 9.3 | 38 |
| Aroclor-1254 | | 16 | U | 16 | 38 |
| Aroclor-1260 | | 10 | U | 10 | 38 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 83 | | 13 - 134 |
| DCB Decachlorobiphenyl | 89 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-11-14**

Lab Sample ID: 240-57597-11 Date Sampled: 11/05/2015 0920
Client Matrix: Solid % Moisture: 16.5 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.28 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1612 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 24 | U | 24 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.5 | U | 9.5 | 39 |
| Aroclor-1254 | | 17 | U | 17 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 82 | | 13 - 134 |
| DCB Decachlorobiphenyl | 95 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-12-14**

Lab Sample ID: 240-57597-12 Date Sampled: 11/05/2015 0922
Client Matrix: Solid % Moisture: 15.9 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.29 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1628 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 24 | U | 24 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.4 | U | 9.4 | 39 |
| Aroclor-1254 | | 16 | U | 16 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 81 | | 13 - 134 |
| DCB Decachlorobiphenyl | 92 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-13-14**

Lab Sample ID: 240-57597-13 Date Sampled: 11/05/2015 0924
Client Matrix: Solid % Moisture: 14.7 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.41 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1644 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 38 |
| Aroclor-1221 | | 19 | U | 19 | 38 |
| Aroclor-1232 | | 23 | U | 23 | 38 |
| Aroclor-1242 | | 13 | U | 13 | 38 |
| Aroclor-1248 | | 9.3 | U | 9.3 | 38 |
| Aroclor-1254 | | 16 | U | 16 | 38 |
| Aroclor-1260 | | 10 | U | 10 | 38 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 77 | | 13 - 134 |
| DCB Decachlorobiphenyl | 94 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-14-14**

Lab Sample ID: 240-57597-14 Date Sampled: 11/05/2015 0926
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.19 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1701 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 38 |
| Aroclor-1221 | | 19 | U | 19 | 38 |
| Aroclor-1232 | | 23 | U | 23 | 38 |
| Aroclor-1242 | | 13 | U | 13 | 38 |
| Aroclor-1248 | | 9.3 | U | 9.3 | 38 |
| Aroclor-1254 | | 16 | U | 16 | 38 |
| Aroclor-1260 | | 10 | U | 10 | 38 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 78 | | 13 - 134 |
| DCB Decachlorobiphenyl | 89 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Client Sample ID:** **F-VB-UST-DUP**

Lab Sample ID: 240-57597-15 Date Sampled: 11/05/2015 0000
Client Matrix: Solid % Moisture: 14.5 Date Received: 11/06/2015 1020

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Prep Method: | 3540C | Prep Batch: | 240-205809 | Initial Weight/Volume: | 30.34 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/11/2015 1806 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/09/2015 0953 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 38 |
| Aroclor-1221 | | 18 | U | 18 | 38 |
| Aroclor-1232 | | 23 | U | 23 | 38 |
| Aroclor-1242 | | 13 | U | 13 | 38 |
| Aroclor-1248 | | 9.2 | U | 9.2 | 38 |
| Aroclor-1254 | | 16 | U | 16 | 38 |
| Aroclor-1260 | | 10 | U | 10 | 38 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 74 | | 13 - 134 |
| DCB Decachlorobiphenyl | 89 | | 10 - 155 |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-1-14**

Lab Sample ID: 240-57597-1 Date Sampled: 11/05/2015 0900
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 84 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 16 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-2-14**

Lab Sample ID: 240-57597-2 Date Sampled: 11/05/2015 0902
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 85 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 15 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-3-14**

Lab Sample ID: 240-57597-3 Date Sampled: 11/05/2015 0904
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 82 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 18 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-4-14**

Lab Sample ID: 240-57597-4 Date Sampled: 11/05/2015 0906
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 84 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 16 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-5-14**

Lab Sample ID: 240-57597-5 Date Sampled: 11/05/2015 0908
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 85 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 15 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-6-14**

Lab Sample ID: 240-57597-6 Date Sampled: 11/05/2015 0910
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 83 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 17 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-7-14**

Lab Sample ID: 240-57597-7 Date Sampled: 11/05/2015 0912
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 84 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 16 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-8-14**

Lab Sample ID: 240-57597-8 Date Sampled: 11/05/2015 0914
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 82 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 18 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-9-14**

Lab Sample ID: 240-57597-9 Date Sampled: 11/05/2015 0916
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 81 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 19 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-10-14**

Lab Sample ID: 240-57597-10 Date Sampled: 11/05/2015 0918
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 87 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 13 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-11-14**

Lab Sample ID: 240-57597-11 Date Sampled: 11/05/2015 0920
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 83 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 17 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-12-14**

Lab Sample ID: 240-57597-12 Date Sampled: 11/05/2015 0922
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 84 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 16 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-13-14**

Lab Sample ID: 240-57597-13 Date Sampled: 11/05/2015 0924
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 85 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 15 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-14-14**

Lab Sample ID: 240-57597-14 Date Sampled: 11/05/2015 0926
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 86 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 14 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-DUP**

Lab Sample ID: 240-57597-15 Date Sampled: 11/05/2015 0000
Client Matrix: Solid Date Received: 11/06/2015 1020

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 86 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |
| Percent Moisture | 14 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-205828 | Analysis Date: 11/09/2015 1019 | | | | | | DryWt Corrected: N |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Surrogate Recovery Report

**8260C Volatile Organic Compounds by GC/MS**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | DBFM %Rec | DCA %Rec | TOL %Rec | BFB %Rec |
|---|---|---|---|---|---|
| 240-57597-1 | F-VB-UST-1-14 | 101 | 103 | 103 | 91 |
| 240-57597-2 | F-VB-UST-2-14 | 95 | 99 | 96 | 85 |
| 240-57597-3 | F-VB-UST-3-14 | 90 | 90 | 93 | 86 |
| 240-57597-4 | F-VB-UST-4-14 | 94 | 102 | 93 | 83 |
| 240-57597-5 | F-VB-UST-5-14 | 101 | 104 | 100 | 87 |
| 240-57597-6 | F-VB-UST-6-14 | 89 | 93 | 88 | 81 |
| 240-57597-7 | F-VB-UST-7-14 | 98 | 103 | 98 | 87 |
| 240-57597-8 | F-VB-UST-8-14 | 93 | 93 | 97 | 85 |
| 240-57597-9 | F-VB-UST-9-14 | 88 | 94 | 89 | 82 |
| 240-57597-10 | F-VB-UST-10-14 | 88 | 93 | 89 | 83 |
| 240-57597-11 | F-VB-UST-11-14 | 87 | 91 | 90 | 81 |
| 240-57597-12 | F-VB-UST-12-14 | 98 | 103 | 99 | 87 |
| 240-57597-13 | F-VB-UST-13-14 | 98 | 100 | 105 | 90 |
| 240-57597-14 | F-VB-UST-14-14 | 89 | 96 | 87 | 78 |
| 240-57597-15 | F-VB-UST-DUP | 101 | 102 | 102 | 89 |
| MB 240-206089/1-A | | 88 | 91 | 91 | 82 |
| MB 240-206267/6 | | 101 | 102 | 101 | 90 |
| LCS 240-206051/5 | | 95 | 92 | 94 | 87 |
| LCS 240-206267/5 | | 99 | 94 | 95 | 89 |
| 240-57597-5 MS | F-VB-UST-5-14 MS | 97 | 92 | 99 | 95 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 97 | 97 | 97 | 88 |
| 240-57597-5 MSD | F-VB-UST-5-14 MSD | 101 | 98 | 101 | 96 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 99 | 95 | 96 | 86 |

| Surrogate | Acceptance Limits |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 37-132 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 58-123 |
| TOL = Toluene-d8 (Surr) | 67-125 |
| BFB = 4-Bromofluorobenzene (Surr) | 52-136 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Surrogate Recovery Report

**8260C Volatile Organic Compounds by GC/MS**

**Client Matrix: Water**

| Lab Sample ID | Client Sample ID | DBFM %Rec | DCA %Rec | TOL %Rec | BFB %Rec |
|---|---|---|---|---|---|
| 240-57597-16 | TB-110515 | 97 | 99 | 85 | 90 |
| MB 240-206461/6 | | 99 | 95 | 88 | 97 |
| LCS 240-206461/4 | | 103 | 99 | 96 | 103 |
| 240-57564-B-1 MS | | 99 | 97 | 88 | 103 |
| 240-57564-B-1 MSD | | 101 | 98 | 90 | 102 |

| Surrogate | Acceptance Limits |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 79-120 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 78-125 |
| TOL = Toluene-d8 (Surr) | 80-120 |
| BFB = 4-Bromofluorobenzene (Surr) | 61-120 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Surrogate Recovery Report

**8270D Semivolatile Organic Compounds (GC/MS)**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | 2FP %Rec | PHL %Rec | NBZ %Rec | FBP %Rec | TBP %Rec | TPH %Rec |
|---|---|---|---|---|---|---|---|
| 240-57597-1 | F-VB-UST-1-14 | 64 | 65 | 62 | 67 | 44 | 76 |
| 240-57597-2 | F-VB-UST-2-14 | 61 | 61 | 58 | 63 | 42 | 78 |
| 240-57597-3 | F-VB-UST-3-14 | 61 | 60 | 58 | 63 | 41 | 73 |
| 240-57597-4 | F-VB-UST-4-14 | 72 | 72 | 69 | 72 | 42 | 85 |
| 240-57597-5 | F-VB-UST-5-14 | 68 | 67 | 64 | 69 | 25 | 77 |
| 240-57597-6 | F-VB-UST-6-14 | 62 | 60 | 62 | 68 | 33 | 79 |
| 240-57597-7 | F-VB-UST-7-14 | 67 | 64 | 62 | 70 | 38 | 80 |
| 240-57597-8 | F-VB-UST-8-14 | 56 | 54 | 52 | 56 | 32 | 70 |
| 240-57597-9 | F-VB-UST-9-14 | 71 | 68 | 68 | 73 | 45 | 87 |
| 240-57597-10 | F-VB-UST-10-14 | 59 | 61 | 60 | 64 | 40 | 78 |
| 240-57597-11 | F-VB-UST-11-14 | 78 | 80 | 80 | 73 | 40 | 88 |
| 240-57597-12 | F-VB-UST-12-14 | 67 | 64 | 67 | 63 | 32 | 72 |
| 240-57597-13 | F-VB-UST-13-14 | 65 | 57 | 65 | 59 | 33 | 77 |
| 240-57597-14 | F-VB-UST-14-14 | 70 | 62 | 73 | 68 | 28 | 79 |
| 240-57597-15 | F-VB-UST-DUP | 71 | 60 | 74 | 68 | 24 | 81 |
| MB 240-205786/23-A | | 66 | 66 | 65 | 68 | 48 | 80 |
| MB 240-205866/8-A | | 81 | 88 | 78 | 76 | 49 | 87 |
| LCS 240-205786/24-A | | 79 | 79 | 79 | 81 | 44 | 93 |
| LCS 240-205866/9-A | | 72 | 75 | 76 | 64 | 33 | 76 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 69 | 62 | 78 | 71 | 30 | 80 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 61 | 53 | 65 | 60 | 36 | 77 |

| Surrogate | Acceptance Limits |
|---|---|
| 2FP = 2-Fluorophenol (Surr) | 24-110 |
| PHL = Phenol-d5 (Surr) | 26-110 |
| NBZ = Nitrobenzene-d5 (Surr) | 20-110 |
| FBP = 2-Fluorobiphenyl (Surr) | 24-110 |
| TBP = 2,4,6-Tribromophenol (Surr) | 10-110 |
| TPH = Terphenyl-d14 (Surr) | 36-110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Surrogate Recovery Report

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | TFT1 %Rec |
|---|---|---|
| 240-57597-1 | F-VB-UST-1-14 | 78 |
| 240-57597-2 | F-VB-UST-2-14 | 79 |
| 240-57597-3 | F-VB-UST-3-14 | 77 |
| 240-57597-4 | F-VB-UST-4-14 | 71 |
| 240-57597-5 | F-VB-UST-5-14 | 71 |
| 240-57597-6 | F-VB-UST-6-14 | 76 |
| 240-57597-7 | F-VB-UST-7-14 | 78 |
| 240-57597-8 | F-VB-UST-8-14 | 73 |
| 240-57597-9 | F-VB-UST-9-14 | 78 |
| 240-57597-10 | F-VB-UST-10-14 | 79 |
| 240-57597-11 | F-VB-UST-11-14 | 79 |
| 240-57597-12 | F-VB-UST-12-14 | 78 |
| 240-57597-13 | F-VB-UST-13-14 | 75 |
| 240-57597-14 | F-VB-UST-14-14 | 80 |
| 240-57597-15 | F-VB-UST-DUP | 74 |
| MB 240-206227/1-A | | 73 |
| MB 240-206444/1-A | | 72 |
| LCS 240-206227/2-A | | 95 |
| LCS 240-206444/2-A | | 100 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 99 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 40 |

| Surrogate | Acceptance Limits |
|---|---|
| TFT = Trifluorotoluene (Surr) | 40-139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Surrogate Recovery Report

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | OTPH1 %Rec |
|---|---|---|
| 240-57597-1 | F-VB-UST-1-14 | 62 |
| 240-57597-2 | F-VB-UST-2-14 | 62 |
| 240-57597-3 | F-VB-UST-3-14 | 59 |
| 240-57597-4 | F-VB-UST-4-14 | 60 |
| 240-57597-5 | F-VB-UST-5-14 | 56 |
| 240-57597-6 | F-VB-UST-6-14 | 60 |
| 240-57597-7 | F-VB-UST-7-14 | 65 |
| 240-57597-8 | F-VB-UST-8-14 | 65 |
| 240-57597-9 | F-VB-UST-9-14 | 57 |
| 240-57597-10 | F-VB-UST-10-14 | 64 |
| 240-57597-11 | F-VB-UST-11-14 | 64 |
| 240-57597-12 | F-VB-UST-12-14 | 73 |
| 240-57597-13 | F-VB-UST-13-14 | 68 |
| 240-57597-14 | F-VB-UST-14-14 | 61 |
| 240-57597-15 | F-VB-UST-DUP | 65 |
| MB 240-205788/22-A | | 56 |
| LCS 240-205788/23-A | | 67 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 81 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 73 |

| Surrogate | Acceptance Limits |
|---|---|
| OTPH = o-Terphenyl (Surr) | 40-160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Surrogate Recovery Report

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | TCX1 %Rec | DCB1 %Rec |
|---|---|---|---|
| 240-57597-1 | F-VB-UST-1-14 | 78 | 87 |
| 240-57597-2 | F-VB-UST-2-14 | 84 | 100 |
| 240-57597-3 | F-VB-UST-3-14 | 68 | 95 |
| 240-57597-4 | F-VB-UST-4-14 | 75 | 104 |
| 240-57597-5 | F-VB-UST-5-14 | 73 | 91 |
| 240-57597-6 | F-VB-UST-6-14 | 76 | 109 |
| 240-57597-7 | F-VB-UST-7-14 | 76 | 98 |
| 240-57597-8 | F-VB-UST-8-14 | 72 | 90 |
| 240-57597-9 | F-VB-UST-9-14 | 77 | 95 |
| 240-57597-10 | F-VB-UST-10-14 | 83 | 89 |
| 240-57597-11 | F-VB-UST-11-14 | 82 | 95 |
| 240-57597-12 | F-VB-UST-12-14 | 81 | 92 |
| 240-57597-13 | F-VB-UST-13-14 | 77 | 94 |
| 240-57597-14 | F-VB-UST-14-14 | 78 | 89 |
| 240-57597-15 | F-VB-UST-DUP | 74 | 89 |
| MB 240-205809/21-A | | 75 | 89 |
| LCS 240-205809/22-A | | 75 | 85 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 81 | 95 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 77 | 89 |

| Surrogate | Acceptance Limits |
|---|---|
| TCX = Tetrachloro-m-xylene | 13-134 |
| DCB = DCB Decachlorobiphenyl | 10-155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Lab Control Sample - Batch: 240-206051**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206051/5 | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146327.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/10/2015 1130 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 20.4 | 102 | 80 - 120 | |
| 1,1,1-Trichloroethane | 20.0 | 22.0 | 110 | 77 - 126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 19.6 | 98 | 77 - 123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 20.9 | 105 | 80 - 138 | |
| 1,1-Dichloroethane | 20.0 | 18.5 | 93 | 76 - 120 | |
| 1,1-Dichloroethene | 20.0 | 19.1 | 95 | 75 - 135 | |
| 1,1-Dichloropropene | 20.0 | 20.3 | 102 | 80 - 120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 19.5 | 98 | 61 - 121 | |
| 1,2,3-Trichloropropane | 20.0 | 20.3 | 102 | 80 - 126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 20.6 | 103 | 64 - 124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 20.3 | 101 | 80 - 129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 19.1 | 95 | 61 - 132 | |
| 1,2-Dichlorobenzene | 20.0 | 19.0 | 95 | 76 - 120 | |
| 1,2-Dichloroethane | 20.0 | 19.3 | 97 | 72 - 120 | |
| 1,2-Dichloropropane | 20.0 | 17.7 | 88 | 80 - 120 | |
| 1,3-Dichlorobenzene | 20.0 | 19.3 | 96 | 78 - 120 | |
| 1,3-Dichloropropane | 20.0 | 18.6 | 93 | 80 - 120 | |
| 1,4-Dichlorobenzene | 20.0 | 18.8 | 94 | 75 - 120 | |
| 2,2-Dichloropropane | 20.0 | 21.3 | 106 | 69 - 135 | |
| 2-Butanone (MEK) | 40.0 | 31.5 | 79 | 52 - 131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 14.8 | 74 | 28 - 173 | J |
| 2-Chlorotoluene | 20.0 | 20.2 | 101 | 78 - 120 | |
| 2-Hexanone | 40.0 | 31.3 | 78 | 64 - 136 | |
| 4-Chlorotoluene | 20.0 | 20.0 | 100 | 79 - 120 | |
| p-Isopropyltoluene | 20.0 | 20.7 | 103 | 77 - 131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 30.7 | 77 | 67 - 135 | |
| Acetone | 40.0 | 32.8 | 82 | 41 - 137 | |
| Benzene | 20.0 | 18.5 | 93 | 79 - 120 | |
| Bromobenzene | 20.0 | 21.3 | 106 | 80 - 120 | |
| Bromoform | 20.0 | 16.3 | 82 | 62 - 133 | |
| Bromomethane | 20.0 | 18.8 | 94 | 42 - 136 | |
| Carbon disulfide | 20.0 | 15.1 | 75 | 62 - 146 | |
| Carbon tetrachloride | 20.0 | 21.9 | 110 | 71 - 129 | |
| Chlorobenzene | 20.0 | 19.9 | 99 | 78 - 120 | |
| Dibromochloromethane | 20.0 | 18.2 | 91 | 72 - 127 | |
| Chloroethane | 20.0 | 18.7 | 93 | 58 - 120 | |
| Chloroform | 20.0 | 19.9 | 100 | 77 - 120 | |
| Chloromethane | 20.0 | 14.5 | 72 | 50 - 120 | |
| cis-1,2-Dichloroethene | 20.0 | 19.6 | 98 | 76 - 120 | |
| cis-1,3-Dichloropropene | 20.0 | 17.3 | 86 | 74 - 128 | |
| Dibromomethane | 20.0 | 19.1 | 96 | 80 - 120 | |
| Bromochloromethane | 20.0 | 21.1 | 106 | 79 - 120 | |
| Dichlorodifluoromethane | 20.0 | 16.4 | 82 | 26 - 120 | |
| Bromodichloromethane | 20.0 | 18.1 | 90 | 80 - 122 | |
| tert-Butyl alcohol (TBA) | 200 | 138 | 69 | 60 - 125 | J |
| Ethylbenzene | 20.0 | 19.8 | 99 | 79 - 120 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Lab Control Sample - Batch: 240-206051**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206051/5 | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146327.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/10/2015 1130 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Hexachlorobutadiene | 20.0 | 21.4 | 107 | 54 - 131 | |
| Isopropylbenzene | 20.0 | 20.6 | 103 | 76 - 122 | |
| Methyl tert-butyl ether | 20.0 | 18.7 | 93 | 49 - 165 | |
| Methylene Chloride | 20.0 | 16.7 | 84 | 75 - 120 | |
| m-Xylene & p-Xylene | 20.0 | 19.8 | 99 | 80 - 120 | |
| Naphthalene | 20.0 | 18.4 | 92 | 65 - 123 | |
| n-Butylbenzene | 20.0 | 20.2 | 101 | 68 - 135 | |
| N-Propylbenzene | 20.0 | 19.8 | 99 | 80 - 129 | |
| o-Xylene | 20.0 | 20.7 | 103 | 80 - 120 | |
| sec-Butylbenzene | 20.0 | 19.9 | 100 | 74 - 129 | |
| Styrene | 20.0 | 18.7 | 93 | 80 - 120 | |
| Tetrachloroethene | 20.0 | 19.6 | 98 | 79 - 120 | |
| Toluene | 20.0 | 20.5 | 102 | 75 - 120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.1 | 100 | 78 - 120 | |
| trans-1,3-Dichloropropene | 20.0 | 19.3 | 97 | 73 - 131 | |
| Trichloroethene | 20.0 | 21.1 | 105 | 79 - 120 | |
| Trichlorofluoromethane | 20.0 | 22.3 | 111 | 57 - 146 | |
| Vinyl acetate | 20.0 | 16.9 | 85 | 70 - 130 | |
| Vinyl chloride | 20.0 | 15.8 | 79 | 57 - 120 | |
| Xylenes, Total | 40.0 | 40.5 | 101 | 80 - 120 | |
| tert-Butylbenzene | 20.0 | 20.1 | 101 | 76 - 126 | |
| 1,2-Dibromoethane | 20.0 | 19.7 | 99 | 80 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 94 | 67 - 125 |
| Dibromofluoromethane (Surr) | 95 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 87 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 92 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-206089**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206089/1-A | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206089 | Lab File ID: | 146331.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.0 g |
| Analysis Date: | 11/10/2015 1256 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.50 | U | 0.50 | 5.0 |
| 1,1,1-Trichloroethane | 0.65 | U | 0.65 | 5.0 |
| 1,1,2,2-Tetrachloroethane | 0.31 | U | 0.31 | 5.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 0.42 | 5.0 |
| 1,1-Dichloroethane | 0.34 | U | 0.34 | 5.0 |
| 1,1-Dichloroethene | 0.80 | U | 0.80 | 5.0 |
| 1,1-Dichloropropene | 0.37 | U | 0.37 | 5.0 |
| 1,2,3-Trichlorobenzene | 0.50 | U | 0.50 | 5.0 |
| 1,2,3-Trichloropropane | 0.78 | U | 0.78 | 5.0 |
| 1,2,3-Trimethylbenzene | 0.46 | U | 0.46 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.38 | U | 0.38 | 5.0 |
| 1,2,4-Trimethylbenzene | 0.58 | U | 0.58 | 5.0 |
| 1,2-Dibromo-3-Chloropropane | 2.7 | U | 2.7 | 10 |
| 1,2-Dichlorobenzene | 0.34 | U | 0.34 | 5.0 |
| 1,2-Dichloroethane | 0.46 | U | 0.46 | 5.0 |
| 1,2-Dichloropropane | 0.14 | U | 0.14 | 5.0 |
| 1,3-Dichlorobenzene | 0.69 | U | 0.69 | 5.0 |
| 1,3-Dichloropropane | 0.40 | U | 0.40 | 5.0 |
| 1,4-Dichlorobenzene | 0.28 | U | 0.28 | 5.0 |
| 2,2-Dichloropropane | 0.44 | U | 0.44 | 5.0 |
| 2-Butanone (MEK) | 2.83 | J | 1.1 | 20 |
| 2-Chloroethyl vinyl ether | 0.64 | U | 0.64 | 50 |
| 2-Chlorotoluene | 0.30 | U | 0.30 | 5.0 |
| 2-Hexanone | 0.68 | U | 0.68 | 20 |
| 4-Chlorotoluene | 0.26 | U | 0.26 | 5.0 |
| p-Isopropyltoluene | 0.41 | U | 0.41 | 5.0 |
| 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 1.2 | 20 |
| Acetone | 5.7 | U | 5.7 | 20 |
| Benzene | 0.90 | U | 0.90 | 5.0 |
| Bromobenzene | 0.33 | U | 0.33 | 5.0 |
| Bromoform | 0.24 | U | 0.24 | 5.0 |
| Bromomethane | 0.42 | U | 0.42 | 5.0 |
| Carbon disulfide | 0.59 | U | 0.59 | 5.0 |
| Carbon tetrachloride | 0.65 | U | 0.65 | 5.0 |
| Chlorobenzene | 0.53 | U | 0.53 | 5.0 |
| Dibromochloromethane | 0.37 | U | 0.37 | 5.0 |
| Chloroethane | 0.44 | U | 0.44 | 5.0 |
| Chloroform | 0.36 | U | 0.36 | 5.0 |
| Chloromethane | 0.79 | U | 0.79 | 5.0 |
| cis-1,2-Dichloroethene | 0.28 | U | 0.28 | 5.0 |
| cis-1,3-Dichloropropene | 0.71 | U | 0.71 | 5.0 |
| Dibromomethane | 0.42 | U | 0.42 | 5.0 |
| Bromochloromethane | 0.33 | U | 0.33 | 5.0 |
| Dichlorodifluoromethane | 0.28 | U | 0.28 | 5.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 5.0 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-206089**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206089/1-A | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206089 | Lab File ID: | 146331.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.0 g |
| Analysis Date: | 11/10/2015 1256 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| tert-Butyl alcohol (TBA) | 5.7 | U | 5.7 | 200 |
| Ethylbenzene | 0.32 | U | 0.32 | 5.0 |
| Hexachlorobutadiene | 0.40 | U | 0.40 | 5.0 |
| Diisopropyl ether | 0.59 | U | 0.59 | 10 |
| Isopropylbenzene | 0.20 | U | 0.20 | 5.0 |
| Methyl tert-butyl ether | 0.43 | U | 0.43 | 5.0 |
| Methylene Chloride | 0.74 | U | 0.74 | 5.0 |
| m-Xylene & p-Xylene | 0.42 | U | 0.42 | 10 |
| Naphthalene | 0.35 | U | 0.35 | 5.0 |
| n-Butylbenzene | 0.59 | U | 0.59 | 5.0 |
| N-Propylbenzene | 0.60 | U | 0.60 | 5.0 |
| o-Xylene | 0.57 | U | 0.57 | 5.0 |
| sec-Butylbenzene | 0.91 | U | 0.91 | 5.0 |
| Styrene | 0.40 | U | 0.40 | 5.0 |
| Tert-amyl methyl ether | 0.43 | U | 0.43 | 5.0 |
| Ethyl t-butyl ether | 0.41 | U | 0.41 | 5.0 |
| Tetrachloroethene | 0.80 | U | 0.80 | 5.0 |
| Toluene | 0.27 | U | 0.27 | 5.0 |
| trans-1,2-Dichloroethene | 0.32 | U | 0.32 | 5.0 |
| trans-1,3-Dichloropropene | 0.29 | U | 0.29 | 5.0 |
| Trichloroethene | 0.38 | U | 0.38 | 5.0 |
| Trichlorofluoromethane | 0.29 | U | 0.29 | 5.0 |
| Vinyl acetate | 0.66 | U | 0.66 | 10 |
| Vinyl chloride | 0.30 | U | 0.30 | 5.0 |
| Xylenes, Total | 0.54 | U | 0.54 | 10 |
| tert-Butylbenzene | 0.67 | U | 0.67 | 5.0 |
| 1,2-Dibromoethane | 0.53 | U | 0.53 | 5.0 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 91 | 67 - 125 |
| Dibromofluoromethane (Surr) | 88 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 82 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 91 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206089**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-5 | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206089 | Lab File ID: | 146341.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.17 g |
| Analysis Date: | 11/10/2015 1630 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-5 | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206089 | Lab File ID: | 146342.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.12 g |
| Analysis Date: | 11/10/2015 1652 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1,1,2-Tetrachloroethane | 101 | 104 | 34 - 135 | 4 | 30 | | |
| 1,1,1-Trichloroethane | 106 | 112 | 51 - 128 | 7 | 30 | | |
| 1,1,2,2-Tetrachloroethane | 117 | 121 | 16 - 179 | 5 | 30 | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 103 | 111 | 50 - 147 | 9 | 30 | | |
| 1,1-Dichloroethane | 93 | 98 | 54 - 122 | 6 | 30 | | |
| 1,1-Dichloroethene | 87 | 95 | 49 - 157 | 10 | 30 | | |
| 1,1-Dichloropropene | 100 | 109 | 50 - 122 | 10 | 30 | | |
| 1,2,3-Trichlorobenzene | 62 | 78 | 10 - 120 | 24 | 30 | | |
| 1,2,3-Trichloropropane | 104 | 125 | 32 - 174 | 19 | 30 | | |
| 1,2,4-Trichlorobenzene | 62 | 86 | 10 - 120 | 33 | 30 | | F2 |
| 1,2,4-Trimethylbenzene | 82 | 109 | 10 - 173 | 29 | 30 | | |
| 1,2-Dibromo-3-Chloropropane | 97 | 117 | 10 - 153 | 19 | 30 | | |
| 1,2-Dichlorobenzene | 79 | 96 | 17 - 122 | 20 | 30 | | |
| 1,2-Dichloroethane | 97 | 105 | 49 - 123 | 9 | 30 | | |
| 1,2-Dichloropropane | 91 | 95 | 61 - 120 | 5 | 30 | | |
| 1,3-Dichlorobenzene | 80 | 98 | 16 - 126 | 21 | 30 | | |
| 1,3-Dichloropropane | 98 | 101 | 54 - 128 | 3 | 30 | | |
| 1,4-Dichlorobenzene | 78 | 97 | 15 - 121 | 22 | 30 | | |
| 2,2-Dichloropropane | 105 | 107 | 49 - 132 | 2 | 30 | | |
| 2-Butanone (MEK) | 71 | 76 | 30 - 143 | 7 | 30 | | |
| 2-Chloroethyl vinyl ether | 72 | 75 | 10 - 162 | 5 | 30 | J | J |
| 2-Chlorotoluene | 91 | 111 | 11 - 162 | 21 | 30 | | |
| 2-Hexanone | 114 | 114 | 37 - 147 | 1 | 30 | | |
| 4-Chlorotoluene | 85 | 108 | 15 - 149 | 25 | 30 | | |
| p-Isopropyltoluene | 83 | 113 | 10 - 165 | 31 | 30 | | F2 |
| 4-Methyl-2-pentanone (MIBK) | 328 | 303 | 43 - 147 | 7 | 30 | F1 | F1 |
| Acetone | 77 | 88 | 24 - 140 | 11 | 30 | | |
| Benzene | 93 | 100 | 53 - 120 | 8 | 30 | | |
| Bromobenzene | 95 | 117 | 24 - 144 | 22 | 30 | | |
| Bromoform | 82 | 86 | 18 - 129 | 7 | 30 | | |
| Bromomethane | 87 | 95 | 33 - 130 | 10 | 30 | | |
| Carbon disulfide | 70 | 76 | 20 - 151 | 10 | 30 | | |
| Carbon tetrachloride | 105 | 115 | 32 - 137 | 10 | 30 | | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206089**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-5 | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206089 | Lab File ID: | 146341.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.17 g |
| Analysis Date: | 11/10/2015 1630 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-5 | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206089 | Lab File ID: | 146342.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.12 g |
| Analysis Date: | 11/10/2015 1652 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Chlorobenzene | 91 | 103 | 37 - 120 | 14 | 30 | | |
| Dibromochloromethane | 89 | 92 | 29 - 135 | 5 | 30 | | |
| Chloroethane | 94 | 97 | 45 - 120 | 4 | 30 | | |
| Chloroform | 100 | 109 | 53 - 120 | 9 | 30 | | |
| Chloromethane | 70 | 74 | 34 - 120 | 8 | 30 | | |
| cis-1,2-Dichloroethene | 97 | 104 | 50 - 120 | 8 | 30 | | |
| cis-1,3-Dichloropropene | 84 | 91 | 27 - 133 | 9 | 30 | | |
| Dibromomethane | 96 | 105 | 52 - 123 | 9 | 30 | | |
| Bromochloromethane | 102 | 113 | 53 - 120 | 11 | 30 | | |
| Dichlorodifluoromethane | 80 | 80 | 17 - 120 | 1 | 30 | | |
| Bromodichloromethane | 88 | 99 | 35 - 132 | 13 | 30 | | |
| tert-Butyl alcohol (TBA) | 72 | 82 | 60 - 125 | 14 | 30 | J | J |
| Ethylbenzene | 90 | 106 | 30 - 131 | 17 | 30 | | |
| Hexachlorobutadiene | 55 | 97 | 10 - 131 | 57 | 30 | | F2 |
| Isopropylbenzene | 86 | 109 | 21 - 134 | 25 | 30 | | |
| Methyl tert-butyl ether | 92 | 97 | 51 - 157 | 6 | 30 | | |
| Methylene Chloride | 82 | 88 | 54 - 120 | 7 | 30 | | |
| m-Xylene & p-Xylene | 85 | 104 | 29 - 131 | 22 | 30 | | |
| Naphthalene | 76 | 89 | 10 - 124 | 16 | 30 | | |
| n-Butylbenzene | 73 | 109 | 10 - 148 | 40 | 30 | | F2 |
| N-Propylbenzene | 86 | 111 | 10 - 178 | 26 | 30 | | |
| o-Xylene | 92 | 109 | 29 - 134 | 18 | 30 | | |
| sec-Butylbenzene | 80 | 109 | 10 - 172 | 32 | 30 | | F2 |
| Styrene | 79 | 95 | 27 - 127 | 20 | 30 | | |
| Tetrachloroethene | 98 | 107 | 31 - 135 | 10 | 30 | | |
| Toluene | 96 | 106 | 39 - 129 | 11 | 30 | | |
| trans-1,2-Dichloroethene | 98 | 107 | 50 - 123 | 10 | 30 | | |
| trans-1,3-Dichloropropene | 93 | 99 | 28 - 137 | 8 | 30 | | |
| Trichloroethene | 101 | 110 | 10 - 177 | 9 | 30 | | |
| Trichlorofluoromethane | 101 | 114 | 36 - 142 | 13 | 30 | | |
| Vinyl acetate | 63 | 77 | 70 - 130 | 21 | 30 | F1 | |
| Vinyl chloride | 76 | 81 | 42 - 120 | 7 | 30 | | |
| Xylenes, Total | 88 | 107 | 30 - 131 | 20 | 30 | | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206089**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-5 | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206089 | Lab File ID: | 146341.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.17 g |
| Analysis Date: | 11/10/2015 1630 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-5 | Analysis Batch: | 240-206051 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206089 | Lab File ID: | 146342.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.12 g |
| Analysis Date: | 11/10/2015 1652 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/10/2015 1124 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| tert-Butylbenzene | 93 | 116 | 10 - 163 | 22 | 30 | | |
| 1,2-Dibromoethane | 106 | 108 | 45 - 127 | 3 | 30 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 99 | 101 | 67 - 125 |
| Dibromofluoromethane (Surr) | 97 | 101 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 95 | 96 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 92 | 98 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206089**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-5 | Units: ug/Kg | MSD Lab Sample ID: | 240-57597-5 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 11/10/2015 1630 | | Analysis Date: | 11/10/2015 1652 | |
| Prep Date: | 11/10/2015 1124 | | Prep Date: | 11/10/2015 1124 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.58 | U | 22.8 | 23.0 | 23.0 | | 24.0 | |
| 1,1,1-Trichloroethane | 0.75 | U | 22.8 | 23.0 | 24.1 | | 25.7 | |
| 1,1,2,2-Tetrachloroethane | 0.36 | U | 22.8 | 23.0 | 26.6 | | 28.0 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 22.8 | 23.0 | 23.5 | | 25.6 | |
| 1,1-Dichloroethane | 0.39 | U | 22.8 | 23.0 | 21.1 | | 22.5 | |
| 1,1-Dichloroethene | 0.93 | U | 22.8 | 23.0 | 19.8 | | 22.0 | |
| 1,1-Dichloropropene | 0.43 | U | 22.8 | 23.0 | 22.7 | | 25.1 | |
| 1,2,3-Trichlorobenzene | 0.58 | U | 22.8 | 23.0 | 14.1 | | 17.9 | |
| 1,2,3-Trichloropropane | 0.90 | U | 22.8 | 23.0 | 23.8 | | 28.9 | |
| 1,2,4-Trichlorobenzene | 0.44 | U | 22.8 | 23.0 | 14.2 | | 19.9 | F2 |
| 1,2,4-Trimethylbenzene | 0.67 | U | 22.8 | 23.0 | 18.7 | | 25.2 | |
| 1,2-Dibromo-3-Chloropropane | 3.1 | U | 22.8 | 23.0 | 22.2 | | 26.9 | |
| 1,2-Dichlorobenzene | 0.39 | U | 22.8 | 23.0 | 18.1 | | 22.2 | |
| 1,2-Dichloroethane | 0.53 | U | 22.8 | 23.0 | 22.0 | | 24.1 | |
| 1,2-Dichloropropane | 0.16 | U | 22.8 | 23.0 | 20.8 | | 21.9 | |
| 1,3-Dichlorobenzene | 0.80 | U | 22.8 | 23.0 | 18.3 | | 22.6 | |
| 1,3-Dichloropropane | 0.46 | U | 22.8 | 23.0 | 22.5 | | 23.2 | |
| 1,4-Dichlorobenzene | 0.32 | U | 22.8 | 23.0 | 17.8 | | 22.3 | |
| 2,2-Dichloropropane | 0.51 | U | 22.8 | 23.0 | 24.0 | | 24.5 | |
| 2-Butanone (MEK) | 4.3 | J | 45.6 | 46.1 | 36.6 | | 39.3 | |
| 2-Chloroethyl vinyl ether | 0.74 | U | 22.8 | 23.0 | 16.4 | J | 17.2 | J |
| 2-Chlorotoluene | 0.35 | U | 22.8 | 23.0 | 20.8 | | 25.6 | |
| 2-Hexanone | 0.79 | U | 45.6 | 46.1 | 52.2 | | 52.6 | |
| 4-Chlorotoluene | 0.30 | U | 22.8 | 23.0 | 19.4 | | 24.9 | |
| p-Isopropyltoluene | 0.47 | U | 22.8 | 23.0 | 19.0 | | 26.0 | F2 |
| 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 45.6 | 46.1 | 150 | F1 | 140 | F1 |
| Acetone | 13 | J | 45.6 | 46.1 | 47.7 | | 53.3 | |
| Benzene | 1.0 | U | 22.8 | 23.0 | 21.1 | | 23.0 | |
| Bromobenzene | 0.38 | U | 22.8 | 23.0 | 21.7 | | 26.9 | |
| Bromoform | 0.28 | U | 22.8 | 23.0 | 18.7 | | 19.9 | |
| Bromomethane | 0.49 | U | 22.8 | 23.0 | 19.8 | | 21.9 | |
| Carbon disulfide | 0.68 | U | 22.8 | 23.0 | 16.0 | | 17.6 | |
| Carbon tetrachloride | 0.75 | U | 22.8 | 23.0 | 24.0 | | 26.5 | |
| Chlorobenzene | 0.61 | U | 22.8 | 23.0 | 20.7 | | 23.7 | |
| Dibromochloromethane | 0.43 | U | 22.8 | 23.0 | 20.2 | | 21.2 | |
| Chloroethane | 0.51 | U | 22.8 | 23.0 | 21.5 | | 22.3 | |
| Chloroform | 0.42 | U | 22.8 | 23.0 | 22.9 | | 25.1 | |
| Chloromethane | 0.92 | U | 22.8 | 23.0 | 15.9 | | 17.1 | |
| cis-1,2-Dichloroethene | 0.32 | U | 22.8 | 23.0 | 22.1 | | 24.0 | |
| cis-1,3-Dichloropropene | 0.82 | U | 22.8 | 23.0 | 19.1 | | 21.0 | |
| Dibromomethane | 0.49 | U | 22.8 | 23.0 | 22.0 | | 24.2 | |
| Bromochloromethane | 0.38 | U | 22.8 | 23.0 | 23.3 | | 26.0 | |
| Dichlorodifluoromethane | 0.32 | U | 22.8 | 23.0 | 18.3 | | 18.5 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206089**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-5 | Units: ug/Kg | MSD Lab Sample ID: | 240-57597-5 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 11/10/2015 1630 | | Analysis Date: | 11/10/2015 1652 | |
| Prep Date: | 11/10/2015 1124 | | Prep Date: | 11/10/2015 1124 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| Bromodichloromethane | 0.34 | U | 22.8 | 23.0 | 20.0 | | 22.9 | |
| tert-Butyl alcohol (TBA) | 6.6 | U | 228 | 230 | 164 | J | 189 | J |
| Ethylbenzene | 0.37 | U | 22.8 | 23.0 | 20.5 | | 24.4 | |
| Hexachlorobutadiene | 0.46 | U | 22.8 | 23.0 | 12.5 | | 22.3 | F2 |
| Isopropylbenzene | 0.23 | U | 22.8 | 23.0 | 19.5 | | 25.1 | |
| Methyl tert-butyl ether | 0.50 | U | 22.8 | 23.0 | 20.9 | | 22.3 | |
| Methylene Chloride | 0.86 | U | 22.8 | 23.0 | 18.8 | | 20.2 | |
| m-Xylene & p-Xylene | 0.49 | U | 22.8 | 23.0 | 19.3 | | 24.0 | |
| Naphthalene | 0.41 | J | 22.8 | 23.0 | 17.8 | | 21.0 | |
| n-Butylbenzene | 0.68 | U | 22.8 | 23.0 | 16.6 | | 25.0 | F2 |
| N-Propylbenzene | 0.70 | U | 22.8 | 23.0 | 19.7 | | 25.5 | |
| o-Xylene | 0.66 | U | 22.8 | 23.0 | 20.9 | | 25.1 | |
| sec-Butylbenzene | 1.1 | U | 22.8 | 23.0 | 18.2 | | 25.1 | F2 |
| Styrene | 0.46 | U | 22.8 | 23.0 | 17.9 | | 21.9 | |
| Tetrachloroethene | 0.93 | U | 22.8 | 23.0 | 22.3 | | 24.7 | |
| Toluene | 0.31 | U | 22.8 | 23.0 | 22.0 | | 24.4 | |
| trans-1,2-Dichloroethene | 0.37 | U | 22.8 | 23.0 | 22.3 | | 24.7 | |
| trans-1,3-Dichloropropene | 0.34 | U | 22.8 | 23.0 | 21.1 | | 22.9 | |
| Trichloroethene | 0.44 | U | 22.8 | 23.0 | 23.2 | | 25.3 | |
| Trichlorofluoromethane | 0.34 | U | 22.8 | 23.0 | 23.0 | | 26.2 | |
| Vinyl acetate | 0.76 | U | 22.8 | 23.0 | 14.4 | F1 | 17.8 | |
| Vinyl chloride | 0.35 | U | 22.8 | 23.0 | 17.4 | | 18.6 | |
| Xylenes, Total | 0.63 | U | 45.6 | 46.1 | 40.2 | | 49.1 | |
| tert-Butylbenzene | 0.78 | U | 22.8 | 23.0 | 21.3 | | 26.7 | |
| 1,2-Dibromoethane | 0.61 | U | 22.8 | 23.0 | 24.2 | | 25.0 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-206267** **Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206267/6 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146361.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/11/2015 1143 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.50 | U | 0.50 | 5.0 |
| 1,1,1-Trichloroethane | 0.65 | U | 0.65 | 5.0 |
| 1,1,2,2-Tetrachloroethane | 0.31 | U | 0.31 | 5.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 0.42 | 5.0 |
| 1,1-Dichloroethane | 0.34 | U | 0.34 | 5.0 |
| 1,1-Dichloroethene | 0.80 | U | 0.80 | 5.0 |
| 1,1-Dichloropropene | 0.37 | U | 0.37 | 5.0 |
| 1,2,3-Trichlorobenzene | 0.50 | U | 0.50 | 5.0 |
| 1,2,3-Trichloropropane | 0.78 | U | 0.78 | 5.0 |
| 1,2,3-Trimethylbenzene | 0.46 | U | 0.46 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.38 | U | 0.38 | 5.0 |
| 1,2,4-Trimethylbenzene | 0.58 | U | 0.58 | 5.0 |
| 1,2-Dibromo-3-Chloropropane | 2.7 | U | 2.7 | 10 |
| 1,2-Dichlorobenzene | 0.34 | U | 0.34 | 5.0 |
| 1,2-Dichloroethane | 0.46 | U | 0.46 | 5.0 |
| 1,2-Dichloropropane | 0.14 | U | 0.14 | 5.0 |
| 1,3-Dichlorobenzene | 0.69 | U | 0.69 | 5.0 |
| 1,3-Dichloropropane | 0.40 | U | 0.40 | 5.0 |
| 1,4-Dichlorobenzene | 0.28 | U | 0.28 | 5.0 |
| 2,2-Dichloropropane | 0.44 | U | 0.44 | 5.0 |
| 2-Butanone (MEK) | 3.14 | J | 1.1 | 20 |
| 2-Chloroethyl vinyl ether | 0.64 | U | 0.64 | 50 |
| 2-Chlorotoluene | 0.30 | U | 0.30 | 5.0 |
| 2-Hexanone | 0.68 | U | 0.68 | 20 |
| 4-Chlorotoluene | 0.26 | U | 0.26 | 5.0 |
| p-Isopropyltoluene | 0.41 | U | 0.41 | 5.0 |
| 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 1.2 | 20 |
| Acetone | 5.7 | U | 5.7 | 20 |
| Benzene | 0.90 | U | 0.90 | 5.0 |
| Bromobenzene | 0.33 | U | 0.33 | 5.0 |
| Bromoform | 0.24 | U | 0.24 | 5.0 |
| Bromomethane | 0.42 | U | 0.42 | 5.0 |
| Carbon disulfide | 0.59 | U | 0.59 | 5.0 |
| Carbon tetrachloride | 0.65 | U | 0.65 | 5.0 |
| Chlorobenzene | 0.53 | U | 0.53 | 5.0 |
| Dibromochloromethane | 0.37 | U | 0.37 | 5.0 |
| Chloroethane | 0.44 | U | 0.44 | 5.0 |
| Chloroform | 0.36 | U | 0.36 | 5.0 |
| Chloromethane | 0.79 | U | 0.79 | 5.0 |
| cis-1,2-Dichloroethene | 0.28 | U | 0.28 | 5.0 |
| cis-1,3-Dichloropropene | 0.71 | U | 0.71 | 5.0 |
| Dibromomethane | 0.42 | U | 0.42 | 5.0 |
| Bromochloromethane | 0.33 | U | 0.33 | 5.0 |
| Dichlorodifluoromethane | 0.28 | U | 0.28 | 5.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 5.0 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-206267**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206267/6 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146361.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/11/2015 1143 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| tert-Butyl alcohol (TBA) | 5.7 | U | 5.7 | 200 |
| Ethylbenzene | 0.32 | U | 0.32 | 5.0 |
| Hexachlorobutadiene | 0.40 | U | 0.40 | 5.0 |
| Diisopropyl ether | 0.59 | U | 0.59 | 10 |
| Isopropylbenzene | 0.20 | U | 0.20 | 5.0 |
| Methyl tert-butyl ether | 0.43 | U | 0.43 | 5.0 |
| Methylene Chloride | 0.74 | U | 0.74 | 5.0 |
| m-Xylene & p-Xylene | 0.42 | U | 0.42 | 10 |
| Naphthalene | 0.350 | J | 0.35 | 5.0 |
| n-Butylbenzene | 0.59 | U | 0.59 | 5.0 |
| N-Propylbenzene | 0.60 | U | 0.60 | 5.0 |
| o-Xylene | 0.57 | U | 0.57 | 5.0 |
| sec-Butylbenzene | 0.91 | U | 0.91 | 5.0 |
| Styrene | 0.40 | U | 0.40 | 5.0 |
| Tert-amyl methyl ether | 0.43 | U | 0.43 | 5.0 |
| Ethyl t-butyl ether | 0.41 | U | 0.41 | 5.0 |
| Tetrachloroethene | 0.80 | U | 0.80 | 5.0 |
| Toluene | 0.369 | J | 0.27 | 5.0 |
| trans-1,2-Dichloroethene | 0.32 | U | 0.32 | 5.0 |
| trans-1,3-Dichloropropene | 0.29 | U | 0.29 | 5.0 |
| Trichloroethene | 0.38 | U | 0.38 | 5.0 |
| Trichlorofluoromethane | 0.29 | U | 0.29 | 5.0 |
| Vinyl acetate | 0.66 | U | 0.66 | 10 |
| Vinyl chloride | 0.30 | U | 0.30 | 5.0 |
| Xylenes, Total | 0.54 | U | 0.54 | 10 |
| tert-Butylbenzene | 0.67 | U | 0.67 | 5.0 |
| 1,2-Dibromoethane | 0.53 | U | 0.53 | 5.0 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 101 | 67 - 125 |
| Dibromofluoromethane (Surr) | 101 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 90 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 102 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Lab Control Sample - Batch: 240-206267**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206267/5 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146360.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/11/2015 1122 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 21.7 | 108 | 80 - 120 | |
| 1,1,1-Trichloroethane | 20.0 | 22.4 | 112 | 77 - 126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 20.2 | 101 | 77 - 123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 21.2 | 106 | 80 - 138 | |
| 1,1-Dichloroethane | 20.0 | 19.0 | 95 | 76 - 120 | |
| 1,1-Dichloroethene | 20.0 | 18.6 | 93 | 75 - 135 | |
| 1,1-Dichloropropene | 20.0 | 20.9 | 104 | 80 - 120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 20.6 | 103 | 61 - 121 | |
| 1,2,3-Trichloropropane | 20.0 | 21.8 | 109 | 80 - 126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 21.3 | 106 | 64 - 124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 21.1 | 106 | 80 - 129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 20.9 | 104 | 61 - 132 | |
| 1,2-Dichlorobenzene | 20.0 | 20.6 | 103 | 76 - 120 | |
| 1,2-Dichloroethane | 20.0 | 20.6 | 103 | 72 - 120 | |
| 1,2-Dichloropropane | 20.0 | 19.1 | 96 | 80 - 120 | |
| 1,3-Dichlorobenzene | 20.0 | 20.2 | 101 | 78 - 120 | |
| 1,3-Dichloropropane | 20.0 | 20.5 | 102 | 80 - 120 | |
| 1,4-Dichlorobenzene | 20.0 | 20.4 | 102 | 75 - 120 | |
| 2,2-Dichloropropane | 20.0 | 21.9 | 110 | 69 - 135 | |
| 2-Butanone (MEK) | 40.0 | 33.7 | 84 | 52 - 131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 15.3 | 76 | 28 - 173 | J |
| 2-Chlorotoluene | 20.0 | 21.5 | 108 | 78 - 120 | |
| 2-Hexanone | 40.0 | 32.9 | 82 | 64 - 136 | |
| 4-Chlorotoluene | 20.0 | 21.6 | 108 | 79 - 120 | |
| p-Isopropyltoluene | 20.0 | 22.3 | 112 | 77 - 131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 32.6 | 82 | 67 - 135 | |
| Acetone | 40.0 | 33.3 | 83 | 41 - 137 | |
| Benzene | 20.0 | 19.4 | 97 | 79 - 120 | |
| Bromobenzene | 20.0 | 22.4 | 112 | 80 - 120 | |
| Bromoform | 20.0 | 17.7 | 89 | 62 - 133 | |
| Bromomethane | 20.0 | 19.6 | 98 | 42 - 136 | |
| Carbon disulfide | 20.0 | 14.8 | 74 | 62 - 146 | |
| Carbon tetrachloride | 20.0 | 22.8 | 114 | 71 - 129 | |
| Chlorobenzene | 20.0 | 21.0 | 105 | 78 - 120 | |
| Dibromochloromethane | 20.0 | 19.4 | 97 | 72 - 127 | |
| Chloroethane | 20.0 | 19.9 | 99 | 58 - 120 | |
| Chloroform | 20.0 | 21.0 | 105 | 77 - 120 | |
| Chloromethane | 20.0 | 15.6 | 78 | 50 - 120 | |
| cis-1,2-Dichloroethene | 20.0 | 20.1 | 101 | 76 - 120 | |
| cis-1,3-Dichloropropene | 20.0 | 18.4 | 92 | 74 - 128 | |
| Dibromomethane | 20.0 | 21.4 | 107 | 80 - 120 | |
| Bromochloromethane | 20.0 | 22.2 | 111 | 79 - 120 | |
| Dichlorodifluoromethane | 20.0 | 20.4 | 102 | 26 - 120 | |
| Bromodichloromethane | 20.0 | 19.2 | 96 | 80 - 122 | |
| tert-Butyl alcohol (TBA) | 200 | 143 | 71 | 60 - 125 | J |
| Ethylbenzene | 20.0 | 21.3 | 106 | 79 - 120 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Lab Control Sample - Batch: 240-206267**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206267/5 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146360.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/11/2015 1122 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Hexachlorobutadiene | 20.0 | 23.6 | 118 | 54 - 131 | |
| Isopropylbenzene | 20.0 | 21.2 | 106 | 76 - 122 | |
| Methyl tert-butyl ether | 20.0 | 18.5 | 93 | 49 - 165 | |
| Methylene Chloride | 20.0 | 16.9 | 84 | 75 - 120 | |
| m-Xylene & p-Xylene | 20.0 | 21.3 | 107 | 80 - 120 | |
| Naphthalene | 20.0 | 18.4 | 92 | 65 - 123 | |
| n-Butylbenzene | 20.0 | 21.8 | 109 | 68 - 135 | |
| N-Propylbenzene | 20.0 | 21.7 | 108 | 80 - 129 | |
| o-Xylene | 20.0 | 21.7 | 108 | 80 - 120 | |
| sec-Butylbenzene | 20.0 | 20.9 | 105 | 74 - 129 | |
| Styrene | 20.0 | 20.0 | 100 | 80 - 120 | |
| Tetrachloroethene | 20.0 | 21.3 | 107 | 79 - 120 | |
| Toluene | 20.0 | 21.3 | 106 | 75 - 120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.3 | 101 | 78 - 120 | |
| trans-1,3-Dichloropropene | 20.0 | 20.6 | 103 | 73 - 131 | |
| Trichloroethene | 20.0 | 21.5 | 108 | 79 - 120 | |
| Trichlorofluoromethane | 20.0 | 22.8 | 114 | 57 - 146 | |
| Vinyl acetate | 20.0 | 17.7 | 88 | 70 - 130 | |
| Vinyl chloride | 20.0 | 16.8 | 84 | 57 - 120 | |
| Xylenes, Total | 40.0 | 43.0 | 108 | 80 - 120 | |
| tert-Butylbenzene | 20.0 | 21.4 | 107 | 76 - 126 | |
| 1,2-Dibromoethane | 20.0 | 21.7 | 109 | 80 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 95 | 67 - 125 |
| Dibromofluoromethane (Surr) | 99 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 89 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 94 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206305**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206305 | Lab File ID: | 146366.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.91 g |
| Analysis Date: | 11/11/2015 1331 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 1103 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206305 | Lab File ID: | 146367.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.17 g |
| Analysis Date: | 11/11/2015 1352 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 1103 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1,1,2-Tetrachloroethane | 102 | 97 | 34 - 135 | 9 | 30 | | |
| 1,1,1-Trichloroethane | 111 | 105 | 51 - 128 | 8 | 30 | | |
| 1,1,2,2-Tetrachloroethane | 97 | 94 | 16 - 179 | 11 | 30 | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 108 | 104 | 50 - 147 | 10 | 30 | | |
| 1,1-Dichloroethane | 93 | 90 | 54 - 122 | 10 | 30 | | |
| 1,1-Dichloroethene | 93 | 90 | 49 - 157 | 10 | 30 | | |
| 1,1-Dichloropropene | 103 | 98 | 50 - 122 | 8 | 30 | | |
| 1,2,3-Trichlorobenzene | 96 | 88 | 10 - 120 | 4 | 30 | | |
| 1,2,3-Trichloropropane | 112 | 91 | 32 - 174 | 8 | 30 | | |
| 1,2,4-Trichlorobenzene | 98 | 92 | 10 - 120 | 6 | 30 | | |
| 1,2,4-Trimethylbenzene | 98 | 92 | 10 - 173 | 7 | 30 | | |
| 1,2-Dibromo-3-Chloropropane | 90 | 91 | 10 - 153 | 14 | 30 | | |
| 1,2-Dichlorobenzene | 94 | 88 | 17 - 122 | 7 | 30 | | |
| 1,2-Dichloroethane | 97 | 93 | 49 - 123 | 9 | 30 | | |
| 1,2-Dichloropropane | 86 | 85 | 61 - 120 | 12 | 30 | | |
| 1,3-Dichlorobenzene | 97 | 84 | 16 - 126 | 1 | 30 | | |
| 1,3-Dichloropropane | 94 | 89 | 54 - 128 | 9 | 30 | | |
| 1,4-Dichlorobenzene | 91 | 85 | 15 - 121 | 7 | 30 | | |
| 2,2-Dichloropropane | 109 | 107 | 49 - 132 | 11 | 30 | | |
| 2-Butanone (MEK) | 67 | 62 | 30 - 143 | 5 | 30 | | |
| 2-Chloroethyl vinyl ether | 63 | 66 | 10 - 162 | 17 | 30 | J | J |
| 2-Chlorotoluene | 101 | 96 | 11 - 162 | 9 | 30 | | |
| 2-Hexanone | 79 | 74 | 37 - 147 | 7 | 30 | | |
| 4-Chlorotoluene | 98 | 89 | 15 - 149 | 4 | 30 | | |
| p-Isopropyltoluene | 102 | 94 | 10 - 165 | 6 | 30 | | |
| 4-Methyl-2-pentanone (MIBK) | 82 | 73 | 43 - 147 | 3 | 30 | | |
| Acetone | 76 | 64 | 24 - 140 | 3 | 30 | | |
| Benzene | 94 | 91 | 53 - 120 | 11 | 30 | | |
| Bromobenzene | 101 | 98 | 24 - 144 | 11 | 30 | | |
| Bromoform | 81 | 77 | 18 - 129 | 8 | 30 | | |
| Bromomethane | 97 | 92 | 33 - 130 | 8 | 30 | | |
| Carbon disulfide | 70 | 68 | 20 - 151 | 11 | 30 | | |
| Carbon tetrachloride | 108 | 105 | 32 - 137 | 10 | 30 | | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206305**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206305 | Lab File ID: | 146366.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.91 g |
| Analysis Date: | 11/11/2015 1331 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 1103 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206305 | Lab File ID: | 146367.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.17 g |
| Analysis Date: | 11/11/2015 1352 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 1103 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| Analyte | % Rec. MS | % Rec. MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
|---|---|---|---|---|---|---|---|
| Chlorobenzene | 97 | 92 | 37 - 120 | 8 | 30 | | |
| Dibromochloromethane | 86 | 86 | 29 - 135 | 13 | 30 | | |
| Chloroethane | 98 | 92 | 45 - 120 | 7 | 30 | | |
| Chloroform | 104 | 100 | 53 - 120 | 9 | 30 | | |
| Chloromethane | 78 | 76 | 34 - 120 | 11 | 30 | | |
| cis-1,2-Dichloroethene | 99 | 95 | 50 - 120 | 9 | 30 | | |
| cis-1,3-Dichloropropene | 81 | 80 | 27 - 133 | 12 | 30 | | |
| Dibromomethane | 108 | 101 | 52 - 123 | 7 | 30 | | |
| Bromochloromethane | 104 | 98 | 53 - 120 | 8 | 30 | | |
| Dichlorodifluoromethane | 101 | 96 | 17 - 120 | 9 | 30 | | |
| Bromodichloromethane | 87 | 90 | 35 - 132 | 17 | 30 | | |
| tert-Butyl alcohol (TBA) | 83 | 70 | 60 - 125 | 4 | 30 | J | J |
| Ethylbenzene | 96 | 93 | 30 - 131 | 10 | 30 | | |
| Hexachlorobutadiene | 106 | 99 | 10 - 131 | 7 | 30 | | |
| Isopropylbenzene | 101 | 96 | 21 - 134 | 8 | 30 | | |
| Methyl tert-butyl ether | 88 | 86 | 51 - 157 | 11 | 30 | | |
| Methylene Chloride | 82 | 85 | 54 - 120 | 17 | 30 | | |
| m-Xylene & p-Xylene | 98 | 89 | 29 - 131 | 4 | 30 | | |
| Naphthalene | 85 | 83 | 10 - 124 | 11 | 30 | | |
| n-Butylbenzene | 98 | 92 | 10 - 148 | 7 | 30 | | |
| N-Propylbenzene | 98 | 91 | 10 - 178 | 7 | 30 | | |
| o-Xylene | 98 | 95 | 29 - 134 | 10 | 30 | | |
| sec-Butylbenzene | 98 | 91 | 10 - 172 | 6 | 30 | | |
| Styrene | 90 | 86 | 27 - 127 | 9 | 30 | | |
| Tetrachloroethene | 109 | 96 | 31 - 135 | 0 | 30 | | |
| Toluene | 102 | 96 | 39 - 129 | 7 | 30 | | |
| trans-1,2-Dichloroethene | 100 | 97 | 50 - 123 | 10 | 30 | | |
| trans-1,3-Dichloropropene | 86 | 87 | 28 - 137 | 14 | 30 | | |
| Trichloroethene | 103 | 99 | 10 - 177 | 10 | 30 | | |
| Trichlorofluoromethane | 107 | 111 | 36 - 142 | 17 | 30 | | |
| Vinyl acetate | 89 | 79 | 70 - 130 | 2 | 30 | | |
| Vinyl chloride | 85 | 82 | 42 - 120 | 9 | 30 | | |
| Xylenes, Total | 98 | 92 | 30 - 131 | 7 | 30 | | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206305**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206305 | Lab File ID: | 146366.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.91 g |
| Analysis Date: | 11/11/2015 1331 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 1103 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206267 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-206305 | Lab File ID: | 146367.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.17 g |
| Analysis Date: | 11/11/2015 1352 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 1103 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| tert-Butylbenzene | 101 | 95 | 10 - 163 | 7 | 30 | | |
| 1,2-Dibromoethane | 100 | 97 | 45 - 127 | 10 | 30 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 97 | 96 | 67 - 125 |
| Dibromofluoromethane (Surr) | 97 | 99 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 88 | 86 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 97 | 95 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206305**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Units: ug/Kg | MSD Lab Sample ID: | 240-57597-15 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 11/11/2015 1331 | | Analysis Date: | 11/11/2015 1352 | |
| Prep Date: | 11/11/2015 1103 | | Prep Date: | 11/11/2015 1103 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.53 | U | 19.8 | 22.6 | 20.2 | | 22.0 | |
| 1,1,1-Trichloroethane | 0.68 | U | 19.8 | 22.6 | 22.0 | | 23.8 | |
| 1,1,2,2-Tetrachloroethane | 0.33 | U | 19.8 | 22.6 | 19.2 | | 21.3 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.44 | U | 19.8 | 22.6 | 21.4 | | 23.5 | |
| 1,1-Dichloroethane | 0.36 | U | 19.8 | 22.6 | 18.4 | | 20.4 | |
| 1,1-Dichloroethene | 0.84 | U | 19.8 | 22.6 | 18.3 | | 20.3 | |
| 1,1-Dichloropropene | 0.39 | U | 19.8 | 22.6 | 20.4 | | 22.2 | |
| 1,2,3-Trichlorobenzene | 0.53 | U | 19.8 | 22.6 | 19.1 | | 19.9 | |
| 1,2,3-Trichloropropane | 0.82 | U | 19.8 | 22.6 | 22.1 | | 20.5 | |
| 1,2,4-Trichlorobenzene | 0.40 | U | 19.8 | 22.6 | 19.5 | | 20.8 | |
| 1,2,4-Trimethylbenzene | 0.61 | U | 19.8 | 22.6 | 19.4 | | 20.8 | |
| 1,2-Dibromo-3-Chloropropane | 2.9 | U | 19.8 | 22.6 | 17.8 | | 20.5 | |
| 1,2-Dichlorobenzene | 0.36 | U | 19.8 | 22.6 | 18.5 | | 19.9 | |
| 1,2-Dichloroethane | 0.48 | U | 19.8 | 22.6 | 19.3 | | 21.1 | |
| 1,2-Dichloropropane | 0.15 | U | 19.8 | 22.6 | 17.0 | | 19.2 | |
| 1,3-Dichlorobenzene | 0.73 | U | 19.8 | 22.6 | 19.2 | | 19.0 | |
| 1,3-Dichloropropane | 0.42 | U | 19.8 | 22.6 | 18.5 | | 20.2 | |
| 1,4-Dichlorobenzene | 0.29 | U | 19.8 | 22.6 | 17.9 | | 19.1 | |
| 2,2-Dichloropropane | 0.46 | U | 19.8 | 22.6 | 21.6 | | 24.1 | |
| 2-Butanone (MEK) | 3.6 | J | 39.6 | 45.2 | 30.1 | | 31.6 | |
| 2-Chloroethyl vinyl ether | 0.67 | U | 19.8 | 22.6 | 12.6 | J | 14.8 | J |
| 2-Chlorotoluene | 0.32 | U | 19.8 | 22.6 | 19.9 | | 21.8 | |
| 2-Hexanone | 0.72 | U | 39.6 | 45.2 | 31.1 | | 33.3 | |
| 4-Chlorotoluene | 0.27 | U | 19.8 | 22.6 | 19.3 | | 20.1 | |
| p-Isopropyltoluene | 0.43 | U | 19.8 | 22.6 | 20.1 | | 21.4 | |
| 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 39.6 | 45.2 | 32.3 | | 33.2 | |
| Acetone | 6.0 | U | 39.6 | 45.2 | 30.0 | | 29.1 | |
| Benzene | 0.95 | U | 19.8 | 22.6 | 18.5 | | 20.6 | |
| Bromobenzene | 0.35 | U | 19.8 | 22.6 | 19.9 | | 22.2 | |
| Bromoform | 0.25 | U | 19.8 | 22.6 | 16.1 | | 17.5 | |
| Bromomethane | 0.44 | U | 19.8 | 22.6 | 19.1 | | 20.8 | |
| Carbon disulfide | 0.62 | U | 19.8 | 22.6 | 13.8 | | 15.4 | |
| Carbon tetrachloride | 0.68 | U | 19.8 | 22.6 | 21.4 | | 23.7 | |
| Chlorobenzene | 0.56 | U | 19.8 | 22.6 | 19.3 | | 20.8 | |
| Dibromochloromethane | 0.39 | U | 19.8 | 22.6 | 17.1 | | 19.4 | |
| Chloroethane | 0.46 | U | 19.8 | 22.6 | 19.4 | | 20.8 | |
| Chloroform | 0.38 | U | 19.8 | 22.6 | 20.5 | | 22.6 | |
| Chloromethane | 0.83 | U | 19.8 | 22.6 | 15.5 | | 17.2 | |
| cis-1,2-Dichloroethene | 0.29 | U | 19.8 | 22.6 | 19.6 | | 21.5 | |
| cis-1,3-Dichloropropene | 0.75 | U | 19.8 | 22.6 | 16.0 | | 18.1 | |
| Dibromomethane | 0.44 | U | 19.8 | 22.6 | 21.4 | | 22.8 | |
| Bromochloromethane | 0.35 | U | 19.8 | 22.6 | 20.6 | | 22.2 | |
| Dichlorodifluoromethane | 0.29 | U | 19.8 | 22.6 | 19.9 | | 21.8 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206305**

**Method: 8260C**
**Preparation: 5030C**

| | | | |
|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Units: ug/Kg | |
| Client Matrix: | Solid | | |
| Dilution: | 1.0 | | |
| Analysis Date: | 11/11/2015 1331 | | |
| Prep Date: | 11/11/2015 1103 | | |
| Leach Date: | N/A | | |

| | |
|---|---|
| MSD Lab Sample ID: | 240-57597-15 |
| Client Matrix: | Solid |
| Dilution: | 1.0 |
| Analysis Date: | 11/11/2015 1352 |
| Prep Date: | 11/11/2015 1103 |
| Leach Date: | N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| Bromodichloromethane | 0.31 | U | 19.8 | 22.6 | 17.2 | | 20.4 | |
| tert-Butyl alcohol (TBA) | 6.0 | U | 198 | 226 | 163 | J | 158 | J |
| Ethylbenzene | 0.34 | U | 19.8 | 22.6 | 19.0 | | 20.9 | |
| Hexachlorobutadiene | 0.42 | U | 19.8 | 22.6 | 20.9 | | 22.4 | |
| Isopropylbenzene | 0.21 | U | 19.8 | 22.6 | 20.0 | | 21.8 | |
| Methyl tert-butyl ether | 0.45 | U | 19.8 | 22.6 | 17.5 | | 19.5 | |
| Methylene Chloride | 0.78 | U | 19.8 | 22.6 | 16.2 | | 19.2 | |
| m-Xylene & p-Xylene | 0.44 | U | 19.8 | 22.6 | 19.4 | | 20.2 | |
| Naphthalene | 0.37 | U | 19.8 | 22.6 | 16.9 | | 18.8 | |
| n-Butylbenzene | 0.62 | U | 19.8 | 22.6 | 19.3 | | 20.7 | |
| N-Propylbenzene | 0.63 | U | 19.8 | 22.6 | 19.4 | | 20.7 | |
| o-Xylene | 0.60 | U | 19.8 | 22.6 | 19.5 | | 21.5 | |
| sec-Butylbenzene | 0.96 | U | 19.8 | 22.6 | 19.5 | | 20.6 | |
| Styrene | 0.42 | U | 19.8 | 22.6 | 17.8 | | 19.5 | |
| Tetrachloroethene | 0.84 | U | 19.8 | 22.6 | 21.7 | | 21.7 | |
| Toluene | 0.28 | U | 19.8 | 22.6 | 20.2 | | 21.7 | |
| trans-1,2-Dichloroethene | 0.34 | U | 19.8 | 22.6 | 19.7 | | 21.8 | |
| trans-1,3-Dichloropropene | 0.31 | U | 19.8 | 22.6 | 17.1 | | 19.7 | |
| Trichloroethene | 0.40 | U | 19.8 | 22.6 | 20.3 | | 22.5 | |
| Trichlorofluoromethane | 0.31 | U | 19.8 | 22.6 | 21.2 | | 25.1 | |
| Vinyl acetate | 0.70 | U | 19.8 | 22.6 | 17.6 | | 17.9 | |
| Vinyl chloride | 0.32 | U | 19.8 | 22.6 | 16.9 | | 18.5 | |
| Xylenes, Total | 0.57 | U | 39.6 | 45.2 | 38.9 | | 41.7 | |
| tert-Butylbenzene | 0.71 | U | 19.8 | 22.6 | 20.0 | | 21.5 | |
| 1,2-Dibromoethane | 0.56 | U | 19.8 | 22.6 | 19.8 | | 21.9 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-206461**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206461/6 | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC4396.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 1211 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 1211 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.28 | U | 0.28 | 1.0 |
| 1,1,1-Trichloroethane | 0.44 | U | 0.44 | 1.0 |
| 1,1,2,2-Tetrachloroethane | 0.22 | U | 0.22 | 1.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloroethane | 0.30 | U | 0.30 | 1.0 |
| 1,1-Dichloroethene | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloropropene | 0.42 | U | 0.42 | 1.0 |
| 1,2,3-Trichlorobenzene | 0.37 | U | 0.37 | 1.0 |
| 1,2,3-Trichloropropane | 0.44 | U | 0.44 | 1.0 |
| 1,2,3-Trimethylbenzene | 0.47 | U | 0.47 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.32 | U | 0.32 | 1.0 |
| 1,2,4-Trimethylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromo-3-Chloropropane | 0.82 | U | 0.82 | 2.0 |
| 1,2-Dichlorobenzene | 0.25 | U | 0.25 | 1.0 |
| 1,2-Dichloroethane | 0.23 | U | 0.23 | 1.0 |
| 1,2-Dichloropropane | 0.25 | U | 0.25 | 1.0 |
| 1,3-Dichlorobenzene | 0.19 | U | 0.19 | 1.0 |
| 1,3-Dichloropropane | 0.19 | U | 0.19 | 1.0 |
| 1,4-Dichlorobenzene | 0.27 | U | 0.27 | 1.0 |
| 2,2-Dichloropropane | 0.26 | U | 0.26 | 1.0 |
| 2-Butanone (MEK) | 0.583 | J | 0.53 | 10 |
| 2-Chloroethyl vinyl ether | 0.59 | U | 0.59 | 10 |
| 2-Chlorotoluene | 0.40 | U | 0.40 | 1.0 |
| 2-Hexanone | 0.48 | U | 0.48 | 10 |
| 4-Chlorotoluene | 0.29 | U | 0.29 | 1.0 |
| p-Isopropyltoluene | 0.43 | U | 0.43 | 1.0 |
| 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 0.99 | 10 |
| Acetone | 0.94 | U | 0.94 | 10 |
| Benzene | 0.35 | U | 0.35 | 1.0 |
| Bromobenzene | 0.35 | U | 0.35 | 1.0 |
| Bromoform | 0.56 | U | 0.56 | 1.0 |
| Bromomethane | 0.44 | U | 0.44 | 1.0 |
| Carbon disulfide | 0.38 | U | 0.38 | 1.0 |
| Carbon tetrachloride | 0.43 | U | 0.43 | 1.0 |
| Chlorobenzene | 0.25 | U | 0.25 | 1.0 |
| Dibromochloromethane | 0.43 | U | 0.43 | 1.0 |
| Chloroethane | 0.32 | U | 0.32 | 1.0 |
| Chloroform | 0.25 | U | 0.25 | 1.0 |
| Chloromethane | 0.44 | U | 0.44 | 1.0 |
| cis-1,2-Dichloroethene | 0.26 | U | 0.26 | 1.0 |
| cis-1,3-Dichloropropene | 0.46 | U | 0.46 | 1.0 |
| Dibromomethane | 0.42 | U | 0.42 | 1.0 |
| Bromochloromethane | 0.50 | U | 0.50 | 1.0 |
| Dichlorodifluoromethane | 0.32 | U | 0.32 | 1.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 1.0 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-206461**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206461/6 | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC4396.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 1211 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 1211 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| tert-Butyl alcohol (TBA) | 4.9 | U | 4.9 | 50 |
| Ethylbenzene | 0.25 | U | 0.25 | 1.0 |
| Hexachlorobutadiene | 0.35 | U | 0.35 | 1.0 |
| Diisopropyl ether | 0.50 | U | 0.50 | 10 |
| Isopropylbenzene | 0.35 | U | 0.35 | 1.0 |
| Methyl tert-butyl ether | 0.20 | U | 0.20 | 1.0 |
| Methylene Chloride | 0.33 | U | 0.33 | 1.0 |
| m-Xylene & p-Xylene | 0.24 | U | 0.24 | 2.0 |
| Naphthalene | 0.45 | U | 0.45 | 1.0 |
| n-Butylbenzene | 0.31 | U | 0.31 | 1.0 |
| N-Propylbenzene | 0.40 | U | 0.40 | 1.0 |
| o-Xylene | 0.25 | U | 0.25 | 1.0 |
| sec-Butylbenzene | 0.48 | U | 0.48 | 1.0 |
| Styrene | 0.45 | U | 0.45 | 1.0 |
| Tert-amyl methyl ether | 0.30 | U | 0.30 | 5.0 |
| Ethyl t-butyl ether | 0.23 | U | 0.23 | 5.0 |
| Tetrachloroethene | 0.31 | U | 0.31 | 1.0 |
| Toluene | 0.23 | U | 0.23 | 1.0 |
| trans-1,2-Dichloroethene | 0.30 | U | 0.30 | 1.0 |
| trans-1,3-Dichloropropene | 0.56 | U | 0.56 | 1.0 |
| Trichloroethene | 0.22 | U | 0.22 | 1.0 |
| Trichlorofluoromethane | 0.49 | U | 0.49 | 1.0 |
| Vinyl acetate | 0.41 | U | 0.41 | 2.0 |
| Vinyl chloride | 0.29 | U | 0.29 | 1.0 |
| Xylenes, Total | 0.52 | U | 0.52 | 2.0 |
| tert-Butylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromoethane | 0.32 | U | 0.32 | 1.0 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 88 | 80 - 120 |
| Dibromofluoromethane (Surr) | 99 | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 97 | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 95 | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Lab Control Sample - Batch: 240-206461**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206461/4 | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC4393.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 1103 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 1103 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 10.0 | 11.8 | 118 | 80 - 120 | |
| 1,1,1-Trichloroethane | 10.0 | 11.1 | 111 | 77 - 123 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 9.38 | 94 | 71 - 123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 10.0 | 9.57 | 96 | 67 - 138 | |
| 1,1-Dichloroethane | 10.0 | 9.43 | 94 | 79 - 125 | |
| 1,1-Dichloroethene | 10.0 | 8.73 | 87 | 76 - 124 | |
| 1,1-Dichloropropene | 10.0 | 10.4 | 104 | 80 - 128 | |
| 1,2,3-Trichlorobenzene | 10.0 | 8.64 | 86 | 66 - 120 | |
| 1,2,3-Trichloropropane | 10.0 | 11.3 | 113 | 72 - 126 | |
| 1,2,4-Trichlorobenzene | 10.0 | 8.66 | 87 | 61 - 120 | |
| 1,2,4-Trimethylbenzene | 10.0 | 9.93 | 99 | 76 - 120 | |
| 1,2-Dibromo-3-Chloropropane | 10.0 | 9.17 | 92 | 50 - 132 | |
| 1,2-Dichlorobenzene | 10.0 | 9.76 | 98 | 79 - 120 | |
| 1,2-Dichloroethane | 10.0 | 10.5 | 105 | 80 - 120 | |
| 1,2-Dichloropropane | 10.0 | 9.07 | 91 | 78 - 124 | |
| 1,3-Dichlorobenzene | 10.0 | 9.90 | 99 | 79 - 120 | |
| 1,3-Dichloropropane | 10.0 | 9.29 | 93 | 80 - 120 | |
| 1,4-Dichlorobenzene | 10.0 | 9.83 | 98 | 79 - 120 | |
| 2,2-Dichloropropane | 10.0 | 10.4 | 104 | 71 - 122 | |
| 2-Butanone (MEK) | 20.0 | 16.3 | 82 | 56 - 138 | |
| 2-Chloroethyl vinyl ether | 10.0 | 8.93 | 89 | 54 - 139 | J |
| 2-Chlorotoluene | 10.0 | 10.4 | 104 | 80 - 120 | |
| 2-Hexanone | 20.0 | 17.6 | 88 | 55 - 141 | |
| 4-Chlorotoluene | 10.0 | 9.89 | 99 | 80 - 120 | |
| p-Isopropyltoluene | 10.0 | 9.42 | 94 | 77 - 120 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 18.1 | 90 | 64 - 135 | |
| Acetone | 20.0 | 13.6 | 68 | 34 - 148 | |
| Benzene | 10.0 | 9.23 | 92 | 80 - 120 | |
| Bromobenzene | 10.0 | 10.7 | 107 | 80 - 120 | |
| Bromoform | 10.0 | 10.7 | 107 | 56 - 122 | |
| Bromomethane | 10.0 | 11.0 | 110 | 38 - 132 | |
| Carbon disulfide | 10.0 | 7.86 | 79 | 65 - 144 | |
| Carbon tetrachloride | 10.0 | 11.6 | 116 | 77 - 131 | |
| Chlorobenzene | 10.0 | 10.2 | 102 | 80 - 120 | |
| Dibromochloromethane | 10.0 | 10.6 | 106 | 74 - 120 | |
| Chloroethane | 10.0 | 9.73 | 97 | 36 - 126 | |
| Chloroform | 10.0 | 10.6 | 106 | 80 - 120 | |
| Chloromethane | 10.0 | 8.46 | 85 | 48 - 133 | |
| cis-1,2-Dichloroethene | 10.0 | 9.74 | 97 | 79 - 120 | |
| cis-1,3-Dichloropropene | 10.0 | 9.52 | 95 | 74 - 126 | |
| Dibromomethane | 10.0 | 10.7 | 107 | 80 - 120 | |
| Bromochloromethane | 10.0 | 10.8 | 108 | 80 - 120 | |
| Dichlorodifluoromethane | 10.0 | 10.5 | 105 | 23 - 136 | |
| Bromodichloromethane | 10.0 | 11.0 | 110 | 80 - 120 | |
| tert-Butyl alcohol (TBA) | 100 | 72.2 | 72 | 38 - 144 | |
| Ethylbenzene | 10.0 | 10.5 | 105 | 80 - 120 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Lab Control Sample - Batch: 240-206461**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206461/4 | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC4393.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 1103 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 1103 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Hexachlorobutadiene | 10.0 | 8.37 | 84 | 62 - 120 | |
| Isopropylbenzene | 10.0 | 10.4 | 104 | 77 - 120 | |
| Methyl tert-butyl ether | 10.0 | 10.1 | 101 | 69 - 121 | |
| Methylene Chloride | 10.0 | 9.34 | 93 | 77 - 129 | |
| m-Xylene & p-Xylene | 10.0 | 10.0 | 100 | 80 - 120 | |
| Naphthalene | 10.0 | 8.44 | 84 | 53 - 121 | |
| n-Butylbenzene | 10.0 | 8.71 | 87 | 74 - 120 | |
| N-Propylbenzene | 10.0 | 10.2 | 102 | 79 - 123 | |
| o-Xylene | 10.0 | 10.4 | 104 | 80 - 120 | |
| sec-Butylbenzene | 10.0 | 9.27 | 93 | 77 - 120 | |
| Styrene | 10.0 | 9.85 | 99 | 76 - 122 | |
| Tetrachloroethene | 10.0 | 10.9 | 109 | 78 - 121 | |
| Toluene | 10.0 | 9.77 | 98 | 80 - 120 | |
| trans-1,2-Dichloroethene | 10.0 | 9.53 | 95 | 80 - 124 | |
| trans-1,3-Dichloropropene | 10.0 | 9.90 | 99 | 75 - 131 | |
| Trichloroethene | 10.0 | 11.1 | 111 | 80 - 121 | |
| Trichlorofluoromethane | 10.0 | 9.70 | 97 | 61 - 133 | |
| Vinyl acetate | 10.0 | 9.77 | 98 | 27 - 160 | |
| Vinyl chloride | 10.0 | 9.07 | 91 | 52 - 121 | |
| Xylenes, Total | 20.0 | 20.4 | 102 | 80 - 120 | |
| tert-Butylbenzene | 10.0 | 9.67 | 97 | 77 - 121 | |
| 1,2-Dibromoethane | 10.0 | 10.7 | 107 | 80 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 96 | 80 - 120 |
| Dibromofluoromethane (Surr) | 103 | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 103 | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 99 | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206461**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57564-B-1 MS | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC4417.D |
| Dilution: | 250 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 2005 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 2005 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57564-B-1 MSD | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC4418.D |
| Dilution: | 250 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 2027 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 2027 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1,1-Trichloroethane | 103 | 105 | 69 - 122 | 2 | 14 | | |
| 1,1,2,2-Tetrachloroethane | 85 | 89 | 61 - 130 | 4 | 18 | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 74 | 87 | 44 - 140 | 15 | 35 | | |
| 1,1-Dichloroethane | 91 | 93 | 73 - 124 | 2 | 14 | | |
| 1,1-Dichloroethene | 81 | 84 | 67 - 124 | 4 | 24 | | |
| 1,2,4-Trichlorobenzene | 84 | 88 | 48 - 120 | 5 | 28 | | |
| 1,2-Dibromo-3-Chloropropane | 92 | 100 | 42 - 130 | 8 | 24 | | |
| 1,2-Dichlorobenzene | 92 | 95 | 67 - 118 | 3 | 15 | | |
| 1,2-Dichloroethane | 103 | 104 | 74 - 125 | 0 | 24 | | |
| 1,2-Dichloropropane | 87 | 93 | 73 - 122 | 7 | 15 | | |
| 1,3-Dichlorobenzene | 90 | 91 | 65 - 120 | 1 | 15 | | |
| 1,4-Dichlorobenzene | 92 | 93 | 66 - 120 | 2 | 16 | | |
| 2-Butanone (MEK) | 109 | 116 | 49 - 132 | 3 | 19 | | |
| 2-Hexanone | 89 | 87 | 49 - 142 | 3 | 27 | | |
| 4-Methyl-2-pentanone (MIBK) | 100 | 102 | 58 - 136 | 2 | 32 | | |
| Acetone | 179 | 182 | 32 - 126 | 1 | 28 | F1 | F1 |
| Benzene | 89 | 90 | 73 - 121 | 1 | 13 | | |
| Bromoform | 100 | 99 | 45 - 121 | 1 | 19 | | |
| Bromomethane | 96 | 96 | 26 - 136 | 0 | 35 | | |
| Carbon disulfide | 75 | 77 | 54 - 144 | 3 | 34 | | |
| Carbon tetrachloride | 105 | 111 | 65 - 129 | 6 | 20 | | |
| Chlorobenzene | 96 | 96 | 72 - 120 | 1 | 15 | | |
| Dibromochloromethane | 98 | 98 | 62 - 122 | 0 | 19 | | |
| Chloroethane | 88 | 87 | 27 - 131 | 1 | 35 | | |
| Chloroform | 102 | 103 | 73 - 121 | 1 | 17 | | |
| Chloromethane | 79 | 82 | 39 - 134 | 4 | 20 | | |
| cis-1,2-Dichloroethene | 95 | 97 | 66 - 124 | 2 | 22 | | |
| cis-1,3-Dichloropropene | 90 | 93 | 60 - 120 | 4 | 21 | | |
| Dichlorodifluoromethane | 75 | 89 | 14 - 137 | 18 | 34 | | |
| Bromodichloromethane | 106 | 110 | 72 - 120 | 4 | 19 | | |
| Ethylbenzene | 97 | 95 | 68 - 121 | 3 | 16 | | |
| Isopropylbenzene | 95 | 96 | 61 - 122 | 1 | 20 | | |
| Methyl tert-butyl ether | 101 | 103 | 61 - 121 | 2 | 12 | | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206461**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57564-B-1 MS | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC4417.D |
| Dilution: | 250 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 2005 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 2005 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57564-B-1 MSD | Analysis Batch: | 240-206461 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC4418.D |
| Dilution: | 250 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/12/2015 2027 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 2027 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Methylene Chloride | 94 | 93 | 70 - 124 | 1 | 14 | | |
| m-Xylene & p-Xylene | 96 | 94 | 66 - 123 | 2 | 15 | | |
| o-Xylene | 97 | 97 | 68 - 121 | 0 | 14 | | |
| Styrene | 93 | 93 | 64 - 126 | 1 | 15 | | |
| Tetrachloroethene | 98 | 100 | 59 - 125 | 2 | 20 | | |
| Toluene | 88 | 91 | 72 - 122 | 2 | 15 | | |
| trans-1,2-Dichloroethene | 94 | 96 | 72 - 125 | 3 | 25 | | |
| trans-1,3-Dichloropropene | 93 | 92 | 58 - 132 | 1 | 22 | | |
| Trichloroethene | 103 | 109 | 61 - 129 | 6 | 14 | | |
| Trichlorofluoromethane | 81 | 90 | 49 - 133 | 10 | 25 | | |
| Vinyl chloride | 81 | 82 | 44 - 122 | 1 | 35 | | |
| Xylenes, Total | 96 | 96 | 67 - 122 | 1 | 14 | | |
| 1,2-Dibromoethane | 97 | 99 | 69 - 125 | 2 | 24 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 88 | 90 | 80 - 120 |
| Dibromofluoromethane (Surr) | 99 | 101 | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 103 | 102 | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 97 | 98 | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206461**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57564-B-1 MS | Units: ug/L | MSD Lab Sample ID: | 240-57564-B-1 MSD | |
| Client Matrix: | Water | | Client Matrix: | Water | |
| Dilution: | 250 | | Dilution: | 250 | |
| Analysis Date: | 11/12/2015 2005 | | Analysis Date: | 11/12/2015 2027 | |
| Prep Date: | 11/12/2015 2005 | | Prep Date: | 11/12/2015 2027 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 110 | U | 2500 | 2500 | 2570 | | 2630 | |
| 1,1,2,2-Tetrachloroethane | 55 | U | 2500 | 2500 | 2130 | | 2220 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 110 | U | 2500 | 2500 | 1850 | | 2160 | |
| 1,1-Dichloroethane | 75 | U | 2500 | 2500 | 2270 | | 2310 | |
| 1,1-Dichloroethene | 110 | U | 2500 | 2500 | 2020 | | 2090 | |
| 1,2,4-Trichlorobenzene | 80 | U | 2500 | 2500 | 2090 | | 2190 | |
| 1,2-Dibromo-3-Chloropropane | 210 | U | 2500 | 2500 | 2310 | | 2510 | |
| 1,2-Dichlorobenzene | 63 | U | 2500 | 2500 | 2300 | | 2370 | |
| 1,2-Dichloroethane | 58 | U | 2500 | 2500 | 2580 | | 2590 | |
| 1,2-Dichloropropane | 63 | U | 2500 | 2500 | 2170 | | 2330 | |
| 1,3-Dichlorobenzene | 48 | U | 2500 | 2500 | 2250 | | 2270 | |
| 1,4-Dichlorobenzene | 68 | U | 2500 | 2500 | 2290 | | 2330 | |
| 2-Butanone (MEK) | 7000 | | 5000 | 5000 | 12400 | | 12800 | |
| 2-Hexanone | 120 | U | 5000 | 5000 | 4460 | | 4330 | |
| 4-Methyl-2-pentanone (MIBK) | 370 | J | 5000 | 5000 | 5380 | | 5460 | |
| Acetone | 9800 | | 5000 | 5000 | 18700 | F1 | 18900 | F1 |
| Benzene | 88 | U | 2500 | 2500 | 2220 | | 2250 | |
| Bromoform | 140 | U | 2500 | 2500 | 2500 | | 2490 | |
| Bromomethane | 110 | U | 2500 | 2500 | 2390 | | 2400 | |
| Carbon disulfide | 95 | U | 2500 | 2500 | 1870 | | 1920 | |
| Carbon tetrachloride | 110 | U | 2500 | 2500 | 2620 | | 2790 | |
| Chlorobenzene | 63 | U | 2500 | 2500 | 2390 | | 2400 | |
| Dibromochloromethane | 110 | U | 2500 | 2500 | 2460 | | 2450 | |
| Chloroethane | 80 | U | 2500 | 2500 | 2200 | | 2180 | |
| Chloroform | 63 | U | 2500 | 2500 | 2560 | | 2580 | |
| Chloromethane | 110 | U | 2500 | 2500 | 1980 | | 2060 | |
| cis-1,2-Dichloroethene | 65 | U | 2500 | 2500 | 2370 | | 2420 | |
| cis-1,3-Dichloropropene | 120 | U | 2500 | 2500 | 2250 | | 2330 | |
| Dichlorodifluoromethane | 80 | U | 2500 | 2500 | 1870 | | 2230 | |
| Bromodichloromethane | 73 | U | 2500 | 2500 | 2650 | | 2760 | |
| Ethylbenzene | 63 | U | 2500 | 2500 | 2440 | | 2370 | |
| Isopropylbenzene | 88 | U | 2500 | 2500 | 2380 | | 2400 | |
| Methyl tert-butyl ether | 50 | U | 2500 | 2500 | 2520 | | 2570 | |
| Methylene Chloride | 83 | U | 2500 | 2500 | 2350 | | 2330 | |
| m-Xylene & p-Xylene | 150 | J | 2500 | 2500 | 2540 | | 2510 | |
| o-Xylene | 72 | J | 2500 | 2500 | 2500 | | 2490 | |
| Styrene | 110 | U | 2500 | 2500 | 2320 | | 2340 | |
| Tetrachloroethene | 78 | U | 2500 | 2500 | 2450 | | 2490 | |
| Toluene | 170 | J | 2500 | 2500 | 2380 | | 2430 | |
| trans-1,2-Dichloroethene | 75 | U | 2500 | 2500 | 2340 | | 2400 | |
| trans-1,3-Dichloropropene | 140 | U | 2500 | 2500 | 2330 | | 2310 | |
| Trichloroethene | 55 | U | 2500 | 2500 | 2580 | | 2730 | |
| Trichlorofluoromethane | 120 | U | 2500 | 2500 | 2020 | | 2240 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206461**

**Method: 8260C**
**Preparation: 5030C**

| | | | |
|---|---|---|---|
| MS Lab Sample ID: | 240-57564-B-1 MS | Units: ug/L | |
| Client Matrix: | Water | | |
| Dilution: | 250 | | |
| Analysis Date: | 11/12/2015 2005 | | |
| Prep Date: | 11/12/2015 2005 | | |
| Leach Date: | N/A | | |

| | |
|---|---|
| MSD Lab Sample ID: | 240-57564-B-1 MSD |
| Client Matrix: | Water |
| Dilution: | 250 |
| Analysis Date: | 11/12/2015 2027 |
| Prep Date: | 11/12/2015 2027 |
| Leach Date: | N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| Vinyl chloride | 73 | U | 2500 | 2500 | 2040 | 2050 |
| Xylenes, Total | 220 | J | 5000 | 5000 | 5040 | 5000 |
| 1,2-Dibromoethane | 80 | U | 2500 | 2500 | 2430 | 2470 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-205786**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-205786/23-A | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-205786 | Lab File ID: | 51111004.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 1027 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1'-Biphenyl | 3.5 | U | 3.5 | 50 |
| bis (2-chloroisopropyl) ether | 9.5 | U | 9.5 | 100 |
| 2,4,5-Trichlorophenol | 25 | U | 25 | 150 |
| 2,4,6-Trichlorophenol | 8.9 | U | 8.9 | 150 |
| 2,4-Dichlorophenol | 20 | U | 20 | 150 |
| 2,4-Dimethylphenol | 20 | U | 20 | 150 |
| 2,4-Dinitrophenol | 21 | U | 21 | 330 |
| 2,4-Dinitrotoluene | 17 | U | 17 | 200 |
| 2,6-Dinitrotoluene | 21 | U | 21 | 200 |
| 2-Chloronaphthalene | 0.45 | U | 0.45 | 50 |
| 2-Chlorophenol | 8.2 | U | 8.2 | 50 |
| 2-Methylnaphthalene | 0.50 | U | 0.50 | 6.7 |
| 2-Methylphenol | 11 | U | 11 | 200 |
| 2-Nitroaniline | 9.1 | U | 9.1 | 200 |
| 2-Nitrophenol | 8.3 | U | 8.3 | 50 |
| 3,3'-Dichlorobenzidine | 18 | U | 18 | 100 |
| 3-Nitroaniline | 16 | U | 16 | 200 |
| 4,6-Dinitro-2-methylphenol | 9.2 | U | 9.2 | 150 |
| 4-Bromophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Chloro-3-methylphenol | 21 | U | 21 | 150 |
| 4-Chloroaniline | 17 | U | 17 | 150 |
| 4-Chlorophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Nitroaniline | 26 | U | 26 | 200 |
| 4-Nitrophenol | 17 | U | 17 | 330 |
| Acenaphthene | 0.76 | U | 0.76 | 6.7 |
| Acenaphthylene | 0.35 | U | 0.35 | 6.7 |
| Acetophenone | 9.2 | U | 9.2 | 100 |
| Anthracene | 0.78 | U | 0.78 | 6.7 |
| Atrazine | 9.1 | U | 9.1 | 200 |
| Benzaldehyde | 12 | U | 12 | 100 |
| Benzo[a]anthracene | 0.63 | U | 0.63 | 6.7 |
| Benzo[a]pyrene | 0.64 | U | 0.64 | 6.7 |
| Benzo[b]fluoranthene | 0.59 | U | 0.59 | 6.7 |
| Benzo[g,h,i]perylene | 0.35 | U | 0.35 | 6.7 |
| Benzo[k]fluoranthene | 0.68 | U | 0.68 | 6.7 |
| Bis(2-chloroethoxy)methane | 22 | U | 22 | 100 |
| Bis(2-chloroethyl)ether | 2.0 | U | 2.0 | 100 |
| Bis(2-ethylhexyl) phthalate | 19 | U | 19 | 70 |
| Butyl benzyl phthalate | 10 | U | 10 | 70 |
| Caprolactam | 37 | U | 37 | 330 |
| Carbazole | 27 | U | 27 | 50 |
| Chrysene | 1.1 | U | 1.1 | 6.7 |
| Dibenz(a,h)anthracene | 0.66 | U | 0.66 | 6.7 |
| Dibenzofuran | 0.66 | U | 0.66 | 50 |
| Diethyl phthalate | 16 | U | 16 | 70 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-205786**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-205786/23-A | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-205786 | Lab File ID: | 51111004.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 1027 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Dimethyl phthalate | 17 | U | 17 | 70 |
| Di-n-butyl phthalate | 15 | U | 15 | 70 |
| Di-n-octyl phthalate | 7.9 | U | 7.9 | 70 |
| Fluoranthene | 0.55 | U | 0.55 | 6.7 |
| Fluorene | 0.53 | U | 0.53 | 6.7 |
| Hexachlorobenzene | 2.1 | U | 2.1 | 6.7 |
| Hexachlorobutadiene | 5.6 | U | 5.6 | 50 |
| Hexachlorocyclopentadiene | 8.1 | U | 8.1 | 330 |
| Hexachloroethane | 9.0 | U | 9.0 | 50 |
| Indeno[1,2,3-cd]pyrene | 0.35 | U | 0.35 | 6.7 |
| Isophorone | 13 | U | 13 | 50 |
| N-Nitrosodi-n-propylamine | 6.3 | U | 6.3 | 50 |
| N-Nitrosodiphenylamine | 21 | U | 21 | 50 |
| Naphthalene | 0.82 | U | 0.82 | 6.7 |
| Nitrobenzene | 2.2 | U | 2.2 | 100 |
| Pentachlorophenol | 9.1 | U | 9.1 | 150 |
| Phenanthrene | 0.73 | U | 0.73 | 6.7 |
| Phenol | 7.3 | U | 7.3 | 50 |
| Pyrene | 0.44 | U | 0.44 | 6.7 |
| 3 & 4 Methylphenol | 20 | U | 20 | 400 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 80 | 36 - 110 |
| Phenol-d5 (Surr) | 66 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 65 | 20 - 110 |
| 2-Fluorophenol (Surr) | 66 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 68 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 48 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Lab Control Sample - Batch: 240-205786**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-205786/24-A | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-205786 | Lab File ID: | 51111005.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 1051 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 525 | 79 | 35 - 110 | |
| bis (2-chloroisopropyl) ether | 667 | 611 | 92 | 29 - 110 | |
| 2,4,5-Trichlorophenol | 667 | 428 | 64 | 25 - 110 | |
| 2,4,6-Trichlorophenol | 667 | 319 | 48 | 12 - 110 | |
| 2,4-Dichlorophenol | 667 | 532 | 80 | 39 - 110 | |
| 2,4-Dimethylphenol | 667 | 462 | 69 | 29 - 110 | |
| 2,4-Dinitrophenol | 1330 | 481 | 36 | 10 - 110 | |
| 2,4-Dinitrotoluene | 667 | 605 | 91 | 48 - 110 | |
| 2,6-Dinitrotoluene | 667 | 563 | 84 | 45 - 110 | |
| 2-Chloronaphthalene | 667 | 523 | 78 | 32 - 110 | |
| 2-Chlorophenol | 667 | 528 | 79 | 37 - 110 | |
| 2-Methylnaphthalene | 667 | 550 | 83 | 36 - 110 | |
| 2-Methylphenol | 667 | 576 | 86 | 41 - 110 | |
| 2-Nitroaniline | 667 | 596 | 89 | 45 - 110 | |
| 2-Nitrophenol | 667 | 584 | 88 | 34 - 110 | |
| 3,3'-Dichlorobenzidine | 1330 | 900 | 68 | 28 - 110 | |
| 3-Nitroaniline | 667 | 497 | 75 | 44 - 110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 716 | 54 | 10 - 110 | |
| 4-Bromophenyl phenyl ether | 667 | 571 | 86 | 39 - 110 | |
| 4-Chloro-3-methylphenol | 667 | 571 | 86 | 48 - 110 | |
| 4-Chloroaniline | 667 | 395 | 59 | 30 - 110 | |
| 4-Chlorophenyl phenyl ether | 667 | 549 | 82 | 40 - 110 | |
| 4-Nitroaniline | 667 | 588 | 88 | 48 - 110 | |
| 4-Nitrophenol | 1330 | 1060 | 80 | 28 - 110 | |
| Acenaphthene | 667 | 533 | 80 | 38 - 110 | |
| Acenaphthylene | 667 | 532 | 80 | 40 - 110 | |
| Acetophenone | 667 | 551 | 83 | 40 - 110 | |
| Anthracene | 667 | 545 | 82 | 48 - 110 | |
| Atrazine | 1330 | 1240 | 93 | 54 - 120 | |
| Benzaldehyde | 1330 | 1150 | 86 | 32 - 110 | |
| Benzo[a]anthracene | 667 | 553 | 83 | 50 - 110 | |
| Benzo[a]pyrene | 667 | 592 | 89 | 44 - 110 | |
| Benzo[b]fluoranthene | 667 | 592 | 89 | 43 - 110 | |
| Benzo[g,h,i]perylene | 667 | 573 | 86 | 51 - 110 | |
| Benzo[k]fluoranthene | 667 | 571 | 86 | 38 - 105 | |
| Bis(2-chloroethoxy)methane | 667 | 537 | 80 | 32 - 110 | |
| Bis(2-chloroethyl)ether | 667 | 533 | 80 | 34 - 110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 662 | 99 | 50 - 110 | |
| Butyl benzyl phthalate | 667 | 604 | 91 | 51 - 110 | |
| Caprolactam | 1330 | 1110 | 83 | 44 - 114 | |
| Carbazole | 667 | 564 | 85 | 50 - 110 | |
| Chrysene | 667 | 550 | 83 | 50 - 110 | |
| Dibenz(a,h)anthracene | 667 | 620 | 93 | 51 - 110 | |
| Dibenzofuran | 667 | 548 | 82 | 43 - 110 | |
| Diethyl phthalate | 667 | 618 | 93 | 52 - 110 | |
| Dimethyl phthalate | 667 | 572 | 86 | 50 - 110 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Lab Control Sample - Batch: 240-205786**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-205786/24-A | Analysis Batch: | 240-206256 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-205786 | Lab File ID: | 51111005.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 1051 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 0612 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | 667 | 583 | 88 | 51 - 110 | |
| Di-n-octyl phthalate | 667 | 628 | 94 | 48 - 110 | |
| Fluoranthene | 667 | 539 | 81 | 51 - 110 | |
| Fluorene | 667 | 550 | 83 | 46 - 110 | |
| Hexachlorobenzene | 667 | 553 | 83 | 43 - 110 | |
| Hexachlorobutadiene | 667 | 559 | 84 | 29 - 110 | |
| Hexachlorocyclopentadiene | 667 | 240 | 36 | 12 - 110 | J |
| Hexachloroethane | 667 | 539 | 81 | 30 - 110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 606 | 91 | 50 - 110 | |
| Isophorone | 667 | 547 | 82 | 36 - 110 | |
| N-Nitrosodi-n-propylamine | 667 | 563 | 84 | 38 - 110 | |
| N-Nitrosodiphenylamine | 1330 | 1180 | 89 | 46 - 110 | |
| Naphthalene | 667 | 537 | 81 | 36 - 110 | |
| Nitrobenzene | 667 | 550 | 83 | 32 - 110 | |
| Pentachlorophenol | 1330 | 584 | 44 | 10 - 110 | |
| Phenanthrene | 667 | 547 | 82 | 49 - 110 | |
| Phenol | 667 | 542 | 81 | 38 - 110 | |
| Pyrene | 667 | 574 | 86 | 49 - 110 | |
| 3 & 4 Methylphenol | 667 | 526 | 79 | 40 - 110 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 93 | 36 - 110 |
| Phenol-d5 (Surr) | 79 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 79 | 20 - 110 |
| 2-Fluorophenol (Surr) | 79 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 81 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 44 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-205866**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-205866/8-A | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-205866 | Lab File ID: | 51111006.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 0954 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1'-Biphenyl | 3.5 | U | 3.5 | 50 |
| bis (2-chloroisopropyl) ether | 9.5 | U | 9.5 | 100 |
| 2,4,5-Trichlorophenol | 25 | U | 25 | 150 |
| 2,4,6-Trichlorophenol | 8.9 | U | 8.9 | 150 |
| 2,4-Dichlorophenol | 20 | U | 20 | 150 |
| 2,4-Dimethylphenol | 20 | U | 20 | 150 |
| 2,4-Dinitrophenol | 21 | U | 21 | 330 |
| 2,4-Dinitrotoluene | 17 | U | 17 | 200 |
| 2,6-Dinitrotoluene | 21 | U | 21 | 200 |
| 2-Chloronaphthalene | 0.45 | U | 0.45 | 50 |
| 2-Chlorophenol | 8.2 | U | 8.2 | 50 |
| 2-Methylnaphthalene | 0.50 | U | 0.50 | 6.7 |
| 2-Methylphenol | 11 | U | 11 | 200 |
| 2-Nitroaniline | 9.1 | U | 9.1 | 200 |
| 2-Nitrophenol | 8.3 | U | 8.3 | 50 |
| 3,3'-Dichlorobenzidine | 18 | U | 18 | 100 |
| 3-Nitroaniline | 16 | U | 16 | 200 |
| 4,6-Dinitro-2-methylphenol | 9.2 | U | 9.2 | 150 |
| 4-Bromophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Chloro-3-methylphenol | 21 | U | 21 | 150 |
| 4-Chloroaniline | 17 | U | 17 | 150 |
| 4-Chlorophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Nitroaniline | 26 | U | 26 | 200 |
| 4-Nitrophenol | 17 | U | 17 | 330 |
| Acenaphthene | 0.76 | U | 0.76 | 6.7 |
| Acenaphthylene | 0.35 | U | 0.35 | 6.7 |
| Acetophenone | 9.2 | U | 9.2 | 100 |
| Anthracene | 0.78 | U | 0.78 | 6.7 |
| Atrazine | 9.1 | U | 9.1 | 200 |
| Benzaldehyde | 12 | U | 12 | 100 |
| Benzo[a]anthracene | 0.63 | U | 0.63 | 6.7 |
| Benzo[a]pyrene | 0.64 | U | 0.64 | 6.7 |
| Benzo[b]fluoranthene | 0.59 | U | 0.59 | 6.7 |
| Benzo[g,h,i]perylene | 0.35 | U | 0.35 | 6.7 |
| Benzo[k]fluoranthene | 0.68 | U | 0.68 | 6.7 |
| Bis(2-chloroethoxy)methane | 22 | U | 22 | 100 |
| Bis(2-chloroethyl)ether | 2.0 | U | 2.0 | 100 |
| Bis(2-ethylhexyl) phthalate | 19 | U | 19 | 70 |
| Butyl benzyl phthalate | 10 | U | 10 | 70 |
| Caprolactam | 37 | U | 37 | 330 |
| Carbazole | 27 | U | 27 | 50 |
| Chrysene | 1.1 | U | 1.1 | 6.7 |
| Dibenz(a,h)anthracene | 0.66 | U | 0.66 | 6.7 |
| Dibenzofuran | 0.66 | U | 0.66 | 50 |
| Diethyl phthalate | 16 | U | 16 | 70 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-205866**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-205866/8-A | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-205866 | Lab File ID: | 51111006.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 0954 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Dimethyl phthalate | 17 | U | 17 | 70 |
| Di-n-butyl phthalate | 15 | U | 15 | 70 |
| Di-n-octyl phthalate | 7.9 | U | 7.9 | 70 |
| Fluoranthene | 0.55 | U | 0.55 | 6.7 |
| Fluorene | 0.53 | U | 0.53 | 6.7 |
| Hexachlorobenzene | 2.1 | U | 2.1 | 6.7 |
| Hexachlorobutadiene | 5.6 | U | 5.6 | 50 |
| Hexachlorocyclopentadiene | 8.1 | U | 8.1 | 330 |
| Hexachloroethane | 9.0 | U | 9.0 | 50 |
| Indeno[1,2,3-cd]pyrene | 0.35 | U | 0.35 | 6.7 |
| Isophorone | 13 | U | 13 | 50 |
| N-Nitrosodi-n-propylamine | 6.3 | U | 6.3 | 50 |
| N-Nitrosodiphenylamine | 21 | U | 21 | 50 |
| Naphthalene | 0.82 | U | 0.82 | 6.7 |
| Nitrobenzene | 2.2 | U | 2.2 | 100 |
| Pentachlorophenol | 9.1 | U | 9.1 | 150 |
| Phenanthrene | 0.73 | U | 0.73 | 6.7 |
| Phenol | 7.3 | U | 7.3 | 50 |
| Pyrene | 0.44 | U | 0.44 | 6.7 |
| 3 & 4 Methylphenol | 20 | U | 20 | 400 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 87 | 36 - 110 |
| Phenol-d5 (Surr) | 88 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 78 | 20 - 110 |
| 2-Fluorophenol (Surr) | 81 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 76 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 49 | 10 - 110 |

**Lab Control Sample - Batch: 240-205866**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-205866/9-A | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-205866 | Lab File ID: | 51111007.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 1019 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 396 | 59 | 35 - 110 | |
| bis (2-chloroisopropyl) ether | 667 | 514 | 77 | 29 - 110 | |
| 2,4,5-Trichlorophenol | 667 | 325 | 49 | 25 - 110 | |
| 2,4,6-Trichlorophenol | 667 | 243 | 37 | 12 - 110 | |
| 2,4-Dichlorophenol | 667 | 446 | 67 | 39 - 110 | |
| 2,4-Dimethylphenol | 667 | 444 | 67 | 29 - 110 | |
| 2,4-Dinitrophenol | 1330 | 464 | 35 | 10 - 110 | |
| 2,4-Dinitrotoluene | 667 | 473 | 71 | 48 - 110 | |
| 2,6-Dinitrotoluene | 667 | 462 | 69 | 45 - 110 | |
| 2-Chloronaphthalene | 667 | 425 | 64 | 32 - 110 | |
| 2-Chlorophenol | 667 | 439 | 66 | 37 - 110 | |
| 2-Methylnaphthalene | 667 | 428 | 64 | 36 - 110 | |
| 2-Methylphenol | 667 | 438 | 66 | 41 - 110 | |
| 2-Nitroaniline | 667 | 548 | 82 | 45 - 110 | |
| 2-Nitrophenol | 667 | 459 | 69 | 34 - 110 | |
| 3,3'-Dichlorobenzidine | 1330 | 946 | 71 | 28 - 110 | |
| 3-Nitroaniline | 667 | 446 | 67 | 44 - 110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 724 | 54 | 10 - 110 | |
| 4-Bromophenyl phenyl ether | 667 | 468 | 70 | 39 - 110 | |
| 4-Chloro-3-methylphenol | 667 | 444 | 67 | 48 - 110 | |
| 4-Chloroaniline | 667 | 366 | 55 | 30 - 110 | |
| 4-Chlorophenyl phenyl ether | 667 | 412 | 62 | 40 - 110 | |
| 4-Nitroaniline | 667 | 452 | 68 | 48 - 110 | |
| 4-Nitrophenol | 1330 | 804 | 60 | 28 - 110 | |
| Acenaphthene | 667 | 386 | 58 | 38 - 110 | |
| Acenaphthylene | 667 | 400 | 60 | 40 - 110 | |
| Acetophenone | 667 | 424 | 64 | 40 - 110 | |
| Anthracene | 667 | 462 | 69 | 48 - 110 | |
| Atrazine | 1330 | 1140 | 86 | 54 - 120 | |
| Benzaldehyde | 1330 | 1090 | 82 | 32 - 110 | |
| Benzo[a]anthracene | 667 | 451 | 68 | 50 - 110 | |
| Benzo[a]pyrene | 667 | 491 | 74 | 44 - 110 | |
| Benzo[b]fluoranthene | 667 | 559 | 84 | 43 - 110 | |
| Benzo[g,h,i]perylene | 667 | 467 | 70 | 51 - 110 | |
| Benzo[k]fluoranthene | 667 | 429 | 64 | 38 - 105 | |
| Bis(2-chloroethoxy)methane | 667 | 489 | 73 | 32 - 110 | |
| Bis(2-chloroethyl)ether | 667 | 503 | 75 | 34 - 110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 544 | 82 | 50 - 110 | |
| Butyl benzyl phthalate | 667 | 512 | 77 | 51 - 110 | |
| Caprolactam | 1330 | 1000 | 75 | 44 - 114 | |
| Carbazole | 667 | 536 | 80 | 50 - 110 | |
| Chrysene | 667 | 442 | 66 | 50 - 110 | |
| Dibenz(a,h)anthracene | 667 | 513 | 77 | 51 - 110 | |
| Dibenzofuran | 667 | 406 | 61 | 43 - 110 | |
| Diethyl phthalate | 667 | 418 | 63 | 52 - 110 | |
| Dimethyl phthalate | 667 | 455 | 68 | 50 - 110 | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Lab Control Sample - Batch: 240-205866**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-205866/9-A | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-205866 | Lab File ID: | 51111007.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 1019 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | 667 | 538 | 81 | 51 - 110 | |
| Di-n-octyl phthalate | 667 | 448 | 67 | 48 - 110 | |
| Fluoranthene | 667 | 500 | 75 | 51 - 110 | |
| Fluorene | 667 | 426 | 64 | 46 - 110 | |
| Hexachlorobenzene | 667 | 496 | 74 | 43 - 110 | |
| Hexachlorobutadiene | 667 | 392 | 59 | 29 - 110 | |
| Hexachlorocyclopentadiene | 667 | 244 | 37 | 12 - 110 | J |
| Hexachloroethane | 667 | 423 | 63 | 30 - 110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 505 | 76 | 50 - 110 | |
| Isophorone | 667 | 513 | 77 | 36 - 110 | |
| N-Nitrosodi-n-propylamine | 667 | 592 | 89 | 38 - 110 | |
| N-Nitrosodiphenylamine | 1330 | 987 | 74 | 46 - 110 | |
| Naphthalene | 667 | 387 | 58 | 36 - 110 | |
| Nitrobenzene | 667 | 506 | 76 | 32 - 110 | |
| Pentachlorophenol | 1330 | 573 | 43 | 10 - 110 | |
| Phenanthrene | 667 | 439 | 66 | 49 - 110 | |
| Phenol | 667 | 537 | 81 | 38 - 110 | |
| Pyrene | 667 | 481 | 72 | 49 - 110 | |
| 3 & 4 Methylphenol | 667 | 432 | 65 | 40 - 110 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 76 | 36 - 110 |
| Phenol-d5 (Surr) | 75 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 76 | 20 - 110 |
| 2-Fluorophenol (Surr) | 72 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 64 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 33 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-205866**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-205866 | Lab File ID: | 51111024.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30.05 g |
| Analysis Date: | 11/11/2015 1652 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-205866 | Lab File ID: | 51111025.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.93 g |
| Analysis Date: | 11/11/2015 1717 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1'-Biphenyl | 66 | 54 | 32 - 110 | 20 | 32 | | |
| bis (2-chloroisopropyl) ether | 83 | 69 | 11 - 110 | 19 | 42 | | |
| 2,4,5-Trichlorophenol | 37 | 27 | 10 - 117 | 33 | 99 | | |
| 2,4,6-Trichlorophenol | 41 | 39 | 10 - 110 | 4 | 38 | | |
| 2,4-Dichlorophenol | 57 | 47 | 10 - 110 | 19 | 34 | | |
| 2,4-Dimethylphenol | 35 | 37 | 10 - 110 | 4 | 31 | | |
| 2,4-Dinitrophenol | 60 | 52 | 10 - 110 | 14 | 99 | | |
| 2,4-Dinitrotoluene | 79 | 75 | 32 - 110 | 4 | 30 | | |
| 2,6-Dinitrotoluene | 77 | 70 | 35 - 110 | 9 | 30 | | |
| 2-Chloronaphthalene | 70 | 58 | 28 - 110 | 18 | 30 | | |
| 2-Chlorophenol | 65 | 54 | 10 - 110 | 18 | 47 | | |
| 2-Methylnaphthalene | 68 | 56 | 10 - 133 | 19 | 42 | | |
| 2-Methylphenol | 50 | 53 | 24 - 110 | 6 | 51 | | |
| 2-Nitroaniline | 91 | 83 | 39 - 110 | 9 | 31 | | |
| 2-Nitrophenol | 76 | 63 | 10 - 110 | 19 | 49 | | |
| 3,3'-Dichlorobenzidine | 76 | 77 | 10 - 110 | 2 | 56 | | |
| 3-Nitroaniline | 68 | 71 | 10 - 110 | 3 | 30 | | |
| 4,6-Dinitro-2-methylphenol | 69 | 64 | 10 - 110 | 8 | 55 | | |
| 4-Bromophenyl phenyl ether | 76 | 68 | 33 - 110 | 11 | 30 | | |
| 4-Chloro-3-methylphenol | 30 | 35 | 25 - 110 | 16 | 54 | | |
| 4-Chloroaniline | 50 | 50 | 10 - 110 | 0 | 36 | | |
| 4-Chlorophenyl phenyl ether | 69 | 61 | 32 - 110 | 12 | 30 | | |
| 4-Nitroaniline | 73 | 69 | 10 - 110 | 5 | 48 | | |
| 4-Nitrophenol | 25 | 29 | 10 - 113 | 15 | 49 | J | |
| Acenaphthene | 64 | 55 | 22 - 110 | 15 | 99 | | |
| Acenaphthylene | 67 | 57 | 24 - 110 | 16 | 99 | | |
| Acetophenone | 70 | 57 | 31 - 110 | 20 | 43 | | |
| Anthracene | 75 | 67 | 20 - 110 | 11 | 99 | | |
| Atrazine | 93 | 83 | 45 - 118 | 11 | 30 | | |
| Benzaldehyde | 87 | 71 | 23 - 110 | 20 | 42 | | |
| Benzo[a]anthracene | 71 | 69 | 10 - 122 | 4 | 99 | | |
| Benzo[a]pyrene | 80 | 72 | 10 - 110 | 10 | 99 | | |
| Benzo[b]fluoranthene | 87 | 82 | 12 - 118 | 5 | 99 | | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-205866**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-205866 | Lab File ID: | 51111024.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30.05 g |
| Analysis Date: | 11/11/2015 1652 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206233 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-205866 | Lab File ID: | 51111025.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.93 g |
| Analysis Date: | 11/11/2015 1717 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/09/2015 1223 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | % Rec. MS | % Rec. MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
|---|---|---|---|---|---|---|---|
| Benzo[g,h,i]perylene | 77 | 70 | 10 - 117 | 9 | 99 | | |
| Benzo[k]fluoranthene | 72 | 62 | 10 - 121 | 14 | 99 | | |
| Bis(2-chloroethoxy)methane | 78 | 64 | 26 - 110 | 21 | 37 | | |
| Bis(2-chloroethyl)ether | 78 | 68 | 21 - 110 | 14 | 55 | | |
| Bis(2-ethylhexyl) phthalate | 87 | 85 | 40 - 110 | 3 | 30 | | |
| Butyl benzyl phthalate | 83 | 79 | 44 - 110 | 5 | 30 | | |
| Caprolactam | 81 | 75 | 10 - 134 | 8 | 32 | | |
| Carbazole | 86 | 77 | 34 - 110 | 10 | 30 | | |
| Chrysene | 70 | 67 | 10 - 125 | 4 | 99 | | |
| Dibenz(a,h)anthracene | 82 | 76 | 14 - 113 | 7 | 99 | | |
| Dibenzofuran | 69 | 59 | 29 - 110 | 15 | 30 | | |
| Diethyl phthalate | 70 | 65 | 42 - 110 | 7 | 30 | | |
| Dimethyl phthalate | 76 | 69 | 41 - 110 | 9 | 30 | | |
| Di-n-butyl phthalate | 88 | 81 | 43 - 110 | 8 | 30 | | |
| Di-n-octyl phthalate | 72 | 66 | 24 - 119 | 8 | 30 | | |
| Fluoranthene | 80 | 74 | 10 - 110 | 7 | 99 | | |
| Fluorene | 70 | 62 | 23 - 110 | 12 | 99 | | |
| Hexachlorobenzene | 80 | 71 | 34 - 110 | 11 | 30 | | |
| Hexachlorobutadiene | 61 | 51 | 25 - 110 | 18 | 34 | | |
| Hexachlorocyclopentadiene | 28 | 24 | 10 - 110 | 16 | 79 | J | J |
| Hexachloroethane | 68 | 53 | 12 - 110 | 25 | 50 | | |
| Indeno[1,2,3-cd]pyrene | 80 | 74 | 10 - 114 | 7 | 99 | | |
| Isophorone | 84 | 69 | 29 - 110 | 19 | 38 | | |
| N-Nitrosodi-n-propylamine | 97 | 79 | 26 - 110 | 21 | 42 | | |
| N-Nitrosodiphenylamine | 81 | 72 | 22 - 110 | 11 | 30 | | |
| Naphthalene | 61 | 51 | 10 - 111 | 17 | 99 | | |
| Nitrobenzene | 81 | 66 | 23 - 110 | 21 | 41 | | |
| Pentachlorophenol | 27 | 39 | 10 - 110 | 37 | 50 | | |
| Phenanthrene | 72 | 65 | 10 - 166 | 9 | 99 | | |
| Phenol | 51 | 53 | 17 - 110 | 6 | 53 | | |
| Pyrene | 77 | 74 | 10 - 147 | 5 | 99 | | |
| 3 & 4 Methylphenol | 47 | 50 | 25 - 110 | 6 | 50 | J | J |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|

Client: Tetra Tech, Inc. Job Number: 240-57597-1

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 80 | 77 | 36 - 110 |
| Phenol-d5 (Surr) | 62 | 53 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 78 | 65 | 20 - 110 |
| 2-Fluorophenol (Surr) | 69 | 61 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 71 | 60 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 30 | 36 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-205866**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Units: ug/Kg | MSD Lab Sample ID: | 240-57597-15 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 11/11/2015 1652 | | Analysis Date: | 11/11/2015 1717 | |
| Prep Date: | 11/09/2015 1223 | | Prep Date: | 11/09/2015 1223 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual |
|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 4.1 | U | 778 | 781 | 511 | | 419 |
| bis (2-chloroisopropyl) ether | 11 | U | 778 | 781 | 646 | | 536 |
| 2,4,5-Trichlorophenol | 29 | U | 778 | 781 | 290 | | 208 |
| 2,4,6-Trichlorophenol | 10 | U | 778 | 781 | 318 | | 307 |
| 2,4-Dichlorophenol | 23 | U | 778 | 781 | 444 | | 368 |
| 2,4-Dimethylphenol | 23 | U | 778 | 781 | 275 | | 286 |
| 2,4-Dinitrophenol | 24 | U | 1560 | 1560 | 940 | | 819 |
| 2,4-Dinitrotoluene | 20 | U | 778 | 781 | 616 | | 589 |
| 2,6-Dinitrotoluene | 24 | U | 778 | 781 | 597 | | 546 |
| 2-Chloronaphthalene | 0.52 | U | 778 | 781 | 546 | | 455 |
| 2-Chlorophenol | 9.6 | U | 778 | 781 | 504 | | 422 |
| 2-Methylnaphthalene | 0.58 | U | 778 | 781 | 528 | | 437 |
| 2-Methylphenol | 13 | U | 778 | 781 | 391 | | 416 |
| 2-Nitroaniline | 11 | U | 778 | 781 | 709 | | 646 |
| 2-Nitrophenol | 9.7 | U | 778 | 781 | 593 | | 492 |
| 3,3'-Dichlorobenzidine | 21 | U | 1560 | 1560 | 1180 | | 1210 |
| 3-Nitroaniline | 19 | U | 778 | 781 | 533 | | 551 |
| 4,6-Dinitro-2-methylphenol | 11 | U | 1560 | 1560 | 1080 | | 996 |
| 4-Bromophenyl phenyl ether | 15 | U | 778 | 781 | 592 | | 532 |
| 4-Chloro-3-methylphenol | 24 | U | 778 | 781 | 231 | | 272 |
| 4-Chloroaniline | 20 | U | 778 | 781 | 391 | | 392 |
| 4-Chlorophenyl phenyl ether | 15 | U | 778 | 781 | 537 | | 476 |
| 4-Nitroaniline | 30 | U | 778 | 781 | 570 | | 541 |
| 4-Nitrophenol | 20 | U | 1560 | 1560 | 383 | J | 446 |
| Acenaphthene | 0.89 | U | 778 | 781 | 501 | | 432 |
| Acenaphthylene | 0.41 | U | 778 | 781 | 518 | | 443 |
| Acetophenone | 11 | U | 778 | 781 | 542 | | 442 |
| Anthracene | 0.91 | U | 778 | 781 | 585 | | 523 |
| Atrazine | 11 | U | 1560 | 1560 | 1450 | | 1300 |
| Benzaldehyde | 14 | U | 1560 | 1560 | 1350 | | 1100 |
| Benzo[a]anthracene | 0.73 | U | 778 | 781 | 556 | | 536 |
| Benzo[a]pyrene | 0.75 | U | 778 | 781 | 623 | | 563 |
| Benzo[b]fluoranthene | 0.69 | U | 778 | 781 | 673 | | 638 |
| Benzo[g,h,i]perylene | 0.41 | U | 778 | 781 | 598 | | 545 |
| Benzo[k]fluoranthene | 0.79 | U | 778 | 781 | 562 | | 486 |
| Bis(2-chloroethoxy)methane | 26 | U | 778 | 781 | 611 | | 497 |
| Bis(2-chloroethyl)ether | 2.3 | U | 778 | 781 | 607 | | 528 |
| Bis(2-ethylhexyl) phthalate | 22 | J | 778 | 781 | 679 | | 661 |
| Butyl benzyl phthalate | 12 | U | 778 | 781 | 649 | | 614 |
| Caprolactam | 43 | U | 1560 | 1560 | 1260 | | 1170 |
| Carbazole | 31 | U | 778 | 781 | 666 | | 604 |
| Chrysene | 1.3 | U | 778 | 781 | 544 | | 521 |
| Dibenz(a,h)anthracene | 0.77 | U | 778 | 781 | 637 | | 596 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-205866**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Units: ug/Kg | MSD Lab Sample ID: | 240-57597-15 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 11/11/2015 1652 | | Analysis Date: | 11/11/2015 1717 | |
| Prep Date: | 11/09/2015 1223 | | Prep Date: | 11/09/2015 1223 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| Dibenzofuran | 0.77 | U | 778 | 781 | 536 | | 463 | |
| Diethyl phthalate | 19 | U | 778 | 781 | 546 | | 507 | |
| Dimethyl phthalate | 20 | U | 778 | 781 | 595 | | 543 | |
| Di-n-butyl phthalate | 17 | U | 778 | 781 | 681 | | 631 | |
| Di-n-octyl phthalate | 9.2 | U | 778 | 781 | 559 | | 517 | |
| Fluoranthene | 0.64 | U | 778 | 781 | 621 | | 577 | |
| Fluorene | 0.62 | U | 778 | 781 | 548 | | 486 | |
| Hexachlorobenzene | 2.4 | U | 778 | 781 | 623 | | 557 | |
| Hexachlorobutadiene | 6.5 | U | 778 | 781 | 472 | | 395 | |
| Hexachlorocyclopentadiene | 9.4 | U | 778 | 781 | 219 | J | 186 | J |
| Hexachloroethane | 10 | U | 778 | 781 | 526 | | 411 | |
| Indeno[1,2,3-cd]pyrene | 0.41 | U | 778 | 781 | 622 | | 578 | |
| Isophorone | 15 | U | 778 | 781 | 651 | | 540 | |
| N-Nitrosodi-n-propylamine | 7.3 | U | 778 | 781 | 757 | | 615 | |
| N-Nitrosodiphenylamine | 24 | U | 1560 | 1560 | 1270 | | 1130 | |
| Naphthalene | 0.96 | U | 778 | 781 | 473 | | 397 | |
| Nitrobenzene | 2.6 | U | 778 | 781 | 633 | | 515 | |
| Pentachlorophenol | 11 | U | 1560 | 1560 | 423 | | 617 | |
| Phenanthrene | 0.85 | U | 778 | 781 | 557 | | 510 | |
| Phenol | 8.5 | U | 778 | 781 | 395 | | 418 | |
| Pyrene | 0.51 | U | 778 | 781 | 603 | | 575 | |
| 3 & 4 Methylphenol | 23 | U | 778 | 781 | 368 | J | 392 | J |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-206227**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206227/1-A | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-206227 | Lab File ID: | A0111107.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/11/2015 0956 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 46 | U | 46 | 100 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 73 | 40 - 139 |

**Lab Control Sample - Batch: 240-206227**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206227/2-A | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-206227 | Lab File ID: | A0111108.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/11/2015 1031 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 786 | 98 | 74 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 95 | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206227**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-206227 | Lab File ID: | A0111126.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.04 g |
| Analysis Date: | 11/11/2015 2140 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206211 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-206227 | Lab File ID: | A0111127.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.05 g |
| Analysis Date: | 11/11/2015 2217 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/11/2015 0658 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Gasoline Range Organics [C6 - C10] | 92 | 22 | 10 - 136 | 122 | 40 | | F2 |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 99 | 40 | 40 - 139 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206227**

**Method: 8015C**
**Preparation: 5030A**

| | | | |
|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Units: | ug/Kg |
| Client Matrix: | Solid | | |
| Dilution: | 1.0 | | |
| Analysis Date: | 11/11/2015 2140 | | |
| Prep Date: | 11/11/2015 0658 | | |
| Leach Date: | N/A | | |

| | |
|---|---|
| MSD Lab Sample ID: | 240-57597-15 |
| Client Matrix: | Solid |
| Dilution: | 1.0 |
| Analysis Date: | 11/11/2015 2217 |
| Prep Date: | 11/11/2015 0658 |
| Leach Date: | N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 54 | U | 928 | 926 | 856 | 206 | F2 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-206444**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206444/1-A | Analysis Batch: | 240-206404 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-206444 | Lab File ID: | A0111205.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/12/2015 0803 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 0559 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 46 | U | 46 | 100 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 72 | 40 - 139 |

**Lab Control Sample - Batch: 240-206444**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206444/2-A | Analysis Batch: | 240-206404 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-206444 | Lab File ID: | A0111206.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/12/2015 0840 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/12/2015 0559 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 787 | 98 | 74 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 100 | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-205788**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-205788/22-A | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-205788 | Lab File ID: | P5B11008.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/10/2015 1659 | Units: | mg/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/09/2015 0940 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | 7.82 | J | 6.4 | 17 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| o-Terphenyl (Surr) | 56 | 40 - 160 |

**Lab Control Sample - Batch: 240-205788**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-205788/23-A | Analysis Batch: | 240-206105 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-205788 | Lab File ID: | P5B11009.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/10/2015 1731 | Units: | mg/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/09/2015 0940 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Diesel | 84.0 | 53.7 | 64 | 52 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| o-Terphenyl (Surr) | 67 | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-205788**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206356 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-205788 | Lab File ID: | P5B11113.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.63 g |
| Analysis Date: | 11/11/2015 2108 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/09/2015 0940 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206356 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-205788 | Lab File ID: | P5B11114.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30.20 g |
| Analysis Date: | 11/11/2015 2139 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/09/2015 0940 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | % Rec. | | | | | |
|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Diesel | 75 | 80 | 10 - 124 | 4 | 40 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 81 | 73 | 40 - 160 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-205788**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Units: | mg/Kg | MSD Lab Sample ID: | 240-57597-15 |
| Client Matrix: | Solid | | | Client Matrix: | Solid |
| Dilution: | 1.0 | | | Dilution: | 1.0 |
| Analysis Date: | 11/11/2015 2108 | | | Analysis Date: | 11/11/2015 2139 |
| Prep Date: | 11/09/2015 0940 | | | Prep Date: | 11/09/2015 0940 |
| Leach Date: | N/A | | | Leach Date: | N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| Diesel | 7.5 | U | 99.5 | 97.6 | 75.1 | 77.9 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Method Blank - Batch: 240-205809**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-205809/21-A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Client Matrix: | Solid | Prep Batch: | 240-205809 | Lab File ID: | P4000026.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 1717 | Units: | ug/Kg | Final Weight/Volume: | 10 mL |
| Prep Date: | 11/09/2015 0953 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Aroclor-1016 | 12 | U | 12 | 33 |
| Aroclor-1221 | 16 | U | 16 | 33 |
| Aroclor-1232 | 20 | U | 20 | 33 |
| Aroclor-1242 | 11 | U | 11 | 33 |
| Aroclor-1248 | 8.0 | U | 8.0 | 33 |
| Aroclor-1254 | 14 | U | 14 | 33 |
| Aroclor-1260 | 9.0 | U | 9.0 | 33 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Tetrachloro-m-xylene | 75 | 13 - 134 |
| DCB Decachlorobiphenyl | 89 | 10 - 155 |

**Lab Control Sample - Batch: 240-205809**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-205809/22-A | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Client Matrix: | Solid | Prep Batch: | 240-205809 | Lab File ID: | P4000027.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/11/2015 1733 | Units: | ug/Kg | Final Weight/Volume: | 10 mL |
| Prep Date: | 11/09/2015 0953 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Aroclor-1016 | 333 | 196 | 59 | 51 - 120 | |
| Aroclor-1260 | 334 | 201 | 60 | 48 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Tetrachloro-m-xylene | 75 | 13 - 134 |
| DCB Decachlorobiphenyl | 85 | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-205809**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Client Matrix: | Solid | Prep Batch: | 240-205809 | Lab File ID: | P4000030.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.65 g |
| Analysis Date: | 11/11/2015 1822 | | | Final Weight/Volume: | 10 mL |
| Prep Date: | 11/09/2015 0953 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57597-15 | Analysis Batch: | 240-206283 | Instrument ID: | A2HP4 |
| Client Matrix: | Solid | Prep Batch: | 240-205809 | Lab File ID: | P4000031.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.78 g |
| Analysis Date: | 11/11/2015 1839 | | | Final Weight/Volume: | 10 mL |
| Prep Date: | 11/09/2015 0953 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Aroclor-1016 | 62 | 58 | 17 - 141 | 9 | 56 | | |
| Aroclor-1260 | 66 | 61 | 19 - 137 | 9 | 55 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 81 | 77 | 13 - 134 |
| DCB Decachlorobiphenyl | 95 | 89 | 10 - 155 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-205809**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57597-15 | Units: | ug/Kg | MSD Lab Sample ID: | 240-57597-15 |
| Client Matrix: | Solid | | | Client Matrix: | Solid |
| Dilution: | 1.0 | | | Dilution: | 1.0 |
| Analysis Date: | 11/11/2015 1822 | | | Analysis Date: | 11/11/2015 1839 |
| Prep Date: | 11/09/2015 0953 | | | Prep Date: | 11/09/2015 0953 |
| Leach Date: | N/A | | | Leach Date: | N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| Aroclor-1016 | 14 | U | 394 | 393 | 244 | 229 |
| Aroclor-1260 | 10 | U | 395 | 393 | 260 | 238 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

**Duplicate - Batch: 240-205828**

**Method: Moisture**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | 240-57597-1 | Analysis Batch: | 240-205828 | Instrument ID: | No Equipment Assigned |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | N/A |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | |
| Analysis Date: | 11/09/2015 1019 | Units: | % | Final Weight/Volume: | |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Sample Result/Qual | Result | RPD | Limit | Qual |
|---|---|---|---|---|---|
| Percent Solids | 84 | 84 | 0.2 | 20 | |
| Percent Moisture | 16 | 16 | 0.9 | 20 | |

**Duplicate - Batch: 240-205828**

**Method: Moisture**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | 240-57597-10 | Analysis Batch: | 240-205828 | Instrument ID: | No Equipment Assigned |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | N/A |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | |
| Analysis Date: | 11/09/2015 1019 | Units: | % | Final Weight/Volume: | |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Sample Result/Qual | Result | RPD | Limit | Qual |
|---|---|---|---|---|---|
| Percent Solids | 87 | 87 | 0 | 20 | |
| Percent Moisture | 13 | 13 | 0.2 | 20 | |

# DATA REPORTING QUALIFIERS

Client: Tetra Tech, Inc. Job Number: 240-57597-1

| Lab Section | Qualifier | Description |
|---|---|---|
| GC/MS VOA | | |
| | B | Compound was found in the blank and sample. |
| | U | Indicates the analyte was analyzed for but not detected. |
| | F1 | MS and/or MSD Recovery is outside acceptance limits. |
| | F2 | MS/MSD RPD exceeds control limits |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| GC/MS Semi VOA | | |
| | U | Indicates the analyte was analyzed for but not detected. |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| GC VOA | | |
| | U | Indicates the analyte was analyzed for but not detected. |
| | F2 | MS/MSD RPD exceeds control limits |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| GC Semi VOA | | |
| | B | Compound was found in the blank and sample. |
| | U | Indicates the analyte was analyzed for but not detected. |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC/MS VOA** | | | | | |
| **Analysis Batch:240-206051** | | | | | |
| LCS 240-206051/5 | Lab Control Sample | T | Solid | 8260C | |
| MB 240-206089/1-A | Method Blank | T | Solid | 8260C | 240-206089 |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-5MS | Matrix Spike | T | Solid | 8260C | 240-206089 |
| 240-57597-5MSD | Matrix Spike Duplicate | T | Solid | 8260C | 240-206089 |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 8260C | 240-206089 |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 8260C | 240-206089 |
| **Prep Batch: 240-206089** | | | | | |
| MB 240-206089/1-A | Method Blank | T | Solid | 5030C | |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 5030C | |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 5030C | |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 5030C | |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 5030C | |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 5030C | |
| 240-57597-5MS | Matrix Spike | T | Solid | 5030C | |
| 240-57597-5MSD | Matrix Spike Duplicate | T | Solid | 5030C | |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 5030C | |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 5030C | |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 5030C | |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 5030C | |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 5030C | |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 5030C | |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 5030C | |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 5030C | |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 5030C | |
| **Analysis Batch:240-206267** | | | | | |
| LCS 240-206267/5 | Lab Control Sample | T | Solid | 8260C | |
| MB 240-206267/6 | Method Blank | T | Solid | 8260C | |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 8260C | 240-206305 |
| 240-57597-15MS | Matrix Spike | T | Solid | 8260C | 240-206305 |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 8260C | 240-206305 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC/MS VOA** | | | | | |
| **Prep Batch: 240-206305** | | | | | |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 5030C | |
| 240-57597-15MS | Matrix Spike | T | Solid | 5030C | |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 5030C | |
| **Analysis Batch:240-206461** | | | | | |
| LCS 240-206461/4 | Lab Control Sample | T | Water | 8260C | |
| MB 240-206461/6 | Method Blank | T | Water | 8260C | |
| 240-57564-B-1 MS | Matrix Spike | T | Water | 8260C | |
| 240-57564-B-1 MSD | Matrix Spike Duplicate | T | Water | 8260C | |
| 240-57597-16 | TB-110515 | T | Water | 8260C | |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC/MS Semi VOA** | | | | | |
| **Prep Batch: 240-205786** | | | | | |
| LCS 240-205786/24-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-205786/23-A | Method Blank | T | Solid | 3540C | |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 3540C | |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 3540C | |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 3540C | |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 3540C | |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 3540C | |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 3540C | |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 3540C | |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 3540C | |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 3540C | |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 3540C | |
| **Prep Batch: 240-205866** | | | | | |
| LCS 240-205866/9-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-205866/8-A | Method Blank | T | Solid | 3540C | |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 3540C | |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 3540C | |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 3540C | |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 3540C | |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 3540C | |
| 240-57597-15MS | Matrix Spike | T | Solid | 3540C | |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 3540C | |
| **Analysis Batch:240-206233** | | | | | |
| LCS 240-205866/9-A | Lab Control Sample | T | Solid | 8270D | 240-205866 |
| MB 240-205866/8-A | Method Blank | T | Solid | 8270D | 240-205866 |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 8270D | 240-205866 |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 8270D | 240-205866 |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 8270D | 240-205866 |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 8270D | 240-205866 |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 8270D | 240-205866 |
| 240-57597-15MS | Matrix Spike | T | Solid | 8270D | 240-205866 |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 8270D | 240-205866 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC/MS Semi VOA** | | | | | |
| **Analysis Batch:240-206256** | | | | | |
| LCS 240-205786/24-A | Lab Control Sample | T | Solid | 8270D | 240-205786 |
| MB 240-205786/23-A | Method Blank | T | Solid | 8270D | 240-205786 |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 8270D | 240-205786 |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 8270D | 240-205786 |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 8270D | 240-205786 |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 8270D | 240-205786 |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 8270D | 240-205786 |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 8270D | 240-205786 |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 8270D | 240-205786 |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 8270D | 240-205786 |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 8270D | 240-205786 |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 8270D | 240-205786 |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC VOA** | | | | | |
| **Analysis Batch:240-206211** | | | | | |
| LCS 240-206227/2-A | Lab Control Sample | T | Solid | 8015C | 240-206227 |
| MB 240-206227/1-A | Method Blank | T | Solid | 8015C | 240-206227 |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 8015C | 240-206227 |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 8015C | 240-206227 |
| 240-57597-15MS | Matrix Spike | T | Solid | 8015C | 240-206227 |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 8015C | 240-206227 |
| **Prep Batch: 240-206227** | | | | | |
| LCS 240-206227/2-A | Lab Control Sample | T | Solid | 5030A | |
| MB 240-206227/1-A | Method Blank | T | Solid | 5030A | |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 5030A | |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 5030A | |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 5030A | |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 5030A | |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 5030A | |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 5030A | |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 5030A | |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 5030A | |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 5030A | |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 5030A | |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 5030A | |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 5030A | |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 5030A | |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 5030A | |
| 240-57597-15MS | Matrix Spike | T | Solid | 5030A | |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 5030A | |
| **Analysis Batch:240-206404** | | | | | |
| LCS 240-206444/2-A | Lab Control Sample | T | Solid | 8015C | 240-206444 |
| MB 240-206444/1-A | Method Blank | T | Solid | 8015C | 240-206444 |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 8015C | 240-206444 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC VOA** | | | | | |
| **Prep Batch: 240-206444** | | | | | |
| LCS 240-206444/2-A | Lab Control Sample | T | Solid | 5030A | |
| MB 240-206444/1-A | Method Blank | T | Solid | 5030A | |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 5030A | |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC Semi VOA** | | | | | |
| **Prep Batch: 240-205788** | | | | | |
| LCS 240-205788/23-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-205788/22-A | Method Blank | T | Solid | 3540C | |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 3540C | |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 3540C | |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 3540C | |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 3540C | |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 3540C | |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 3540C | |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 3540C | |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 3540C | |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 3540C | |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 3540C | |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 3540C | |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 3540C | |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 3540C | |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 3540C | |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 3540C | |
| 240-57597-15MS | Matrix Spike | T | Solid | 3540C | |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 3540C | |
| **Prep Batch: 240-205809** | | | | | |
| LCS 240-205809/22-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-205809/21-A | Method Blank | T | Solid | 3540C | |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 3540C | |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 3540C | |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 3540C | |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 3540C | |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 3540C | |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 3540C | |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 3540C | |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 3540C | |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 3540C | |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 3540C | |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 3540C | |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 3540C | |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 3540C | |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 3540C | |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 3540C | |
| 240-57597-15MS | Matrix Spike | T | Solid | 3540C | |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 3540C | |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC Semi VOA** | | | | | |
| **Analysis Batch:240-206105** | | | | | |
| LCS 240-205788/23-A | Lab Control Sample | T | Solid | 8015C | 240-205788 |
| MB 240-205788/22-A | Method Blank | T | Solid | 8015C | 240-205788 |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 8015C | 240-205788 |
| **Analysis Batch:240-206283** | | | | | |
| LCS 240-205809/22-A | Lab Control Sample | T | Solid | 8082A | 240-205809 |
| MB 240-205809/21-A | Method Blank | T | Solid | 8082A | 240-205809 |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 8082A | 240-205809 |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 8082A | 240-205809 |
| 240-57597-15MS | Matrix Spike | T | Solid | 8082A | 240-205809 |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 8082A | 240-205809 |
| **Analysis Batch:240-206356** | | | | | |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | 8015C | 240-205788 |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | 8015C | 240-205788 |
| 240-57597-15MS | Matrix Spike | T | Solid | 8015C | 240-205788 |
| 240-57597-15MSD | Matrix Spike Duplicate | T | Solid | 8015C | 240-205788 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **Report Basis** | | | | | |
| T = Total | | | | | |
| **General Chemistry** | | | | | |
| **Analysis Batch:240-205828** | | | | | |
| 240-57597-1 | F-VB-UST-1-14 | T | Solid | Moisture | |
| 240-57597-1DU | Duplicate | T | Solid | Moisture | |
| 240-57597-2 | F-VB-UST-2-14 | T | Solid | Moisture | |
| 240-57597-3 | F-VB-UST-3-14 | T | Solid | Moisture | |
| 240-57597-4 | F-VB-UST-4-14 | T | Solid | Moisture | |
| 240-57597-5 | F-VB-UST-5-14 | T | Solid | Moisture | |
| 240-57597-6 | F-VB-UST-6-14 | T | Solid | Moisture | |
| 240-57597-7 | F-VB-UST-7-14 | T | Solid | Moisture | |
| 240-57597-8 | F-VB-UST-8-14 | T | Solid | Moisture | |
| 240-57597-9 | F-VB-UST-9-14 | T | Solid | Moisture | |
| 240-57597-10 | F-VB-UST-10-14 | T | Solid | Moisture | |
| 240-57597-10DU | Duplicate | T | Solid | Moisture | |
| 240-57597-11 | F-VB-UST-11-14 | T | Solid | Moisture | |
| 240-57597-12 | F-VB-UST-12-14 | T | Solid | Moisture | |
| 240-57597-13 | F-VB-UST-13-14 | T | Solid | Moisture | |
| 240-57597-14 | F-VB-UST-14-14 | T | Solid | Moisture | |
| 240-57597-15 | F-VB-UST-DUP | T | Solid | Moisture | |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-1** **Client ID: F-VB-UST-1-14**

Sample Date/Time: 11/05/2015 09:00 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-1-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-1-A | | 240-206051 | 240-206089 | 11/10/2015 14:43 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-1-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-1-A | | 240-206256 | 240-205786 | 11/11/2015 11:16 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-1-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-1-A | | 240-206211 | 240-206227 | 11/11/2015 11:08 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-1-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-1-A | | 240-206105 | 240-205788 | 11/10/2015 20:07 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-1-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-1-B | | 240-206283 | 240-205809 | 11/11/2015 13:28 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-1 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-1 DU** **Client ID: F-VB-UST-1-14**

Sample Date/Time: 11/05/2015 09:00 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| A:Moisture | 240-57597-A-1 DU | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-2** **Client ID: F-VB-UST-2-14**

Sample Date/Time: 11/05/2015 09:02 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-2-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-2-A | | 240-206051 | 240-206089 | 11/10/2015 15:05 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-D-2-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-D-2-A | | 240-206256 | 240-205786 | 11/11/2015 11:40 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-2-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-2-A | | 240-206211 | 240-206227 | 11/11/2015 11:45 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-E-2-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-E-2-A | | 240-206105 | 240-205788 | 11/10/2015 20:38 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-D-2-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-D-2-B | | 240-206283 | 240-205809 | 11/11/2015 13:45 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-2 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-3** **Client ID: F-VB-UST-3-14**

Sample Date/Time: 11/05/2015 09:04 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-3-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-3-A | | 240-206051 | 240-206089 | 11/10/2015 15:26 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-3-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-3-A | | 240-206256 | 240-205786 | 11/11/2015 12:05 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-3-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-3-A | | 240-206211 | 240-206227 | 11/11/2015 12:23 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-3-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-3-A | | 240-206105 | 240-205788 | 11/10/2015 21:10 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-3-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-3-B | | 240-206283 | 240-205809 | 11/11/2015 14:01 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-3 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-4** **Client ID: F-VB-UST-4-14**

Sample Date/Time: 11/05/2015 09:06 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-4-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-4-A | | 240-206051 | 240-206089 | 11/10/2015 15:48 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-4-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-4-A | | 240-206256 | 240-205786 | 11/11/2015 12:30 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-4-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-4-A | | 240-206211 | 240-206227 | 11/11/2015 13:00 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-4-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-4-A | | 240-206105 | 240-205788 | 11/10/2015 21:41 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-4-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-4-B | | 240-206283 | 240-205809 | 11/11/2015 14:17 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-4 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-5** **Client ID: F-VB-UST-5-14**

Sample Date/Time: 11/05/2015 09:08 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-5-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-5-A | | 240-206051 | 240-206089 | 11/10/2015 16:09 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-5-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-5-A | | 240-206256 | 240-205786 | 11/11/2015 12:54 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-5-B | | 240-206404 | 240-206444 | 11/12/2015 05:59 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-5-B | | 240-206404 | 240-206444 | 11/12/2015 09:17 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-5-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-5-A | | 240-206105 | 240-205788 | 11/10/2015 22:43 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-5-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-5-B | | 240-206283 | 240-205809 | 11/11/2015 14:33 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-5 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-5 MS** **Client ID: F-VB-UST-5-14**

Sample Date/Time: 11/05/2015 09:08 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-5-B MS | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-5-B MS | | 240-206051 | 240-206089 | 11/10/2015 16:30 | 1 | TAL CAN | TJL2 |

**Lab ID: 240-57597-5 MSD** **Client ID: F-VB-UST-5-14**

Sample Date/Time: 11/05/2015 09:08 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-5-C MSD | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-5-C MSD | | 240-206051 | 240-206089 | 11/10/2015 16:52 | 1 | TAL CAN | TJL2 |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-6** **Client ID: F-VB-UST-6-14**

Sample Date/Time: 11/05/2015 09:10 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-6-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-6-A | | 240-206051 | 240-206089 | 11/10/2015 17:13 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-6-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-6-A | | 240-206256 | 240-205786 | 11/11/2015 13:19 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-6-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-6-A | | 240-206211 | 240-206227 | 11/11/2015 14:13 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-6-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-6-A | | 240-206105 | 240-205788 | 11/10/2015 23:14 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-6-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-6-B | | 240-206283 | 240-205809 | 11/11/2015 14:50 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-6 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-7** **Client ID: F-VB-UST-7-14**

Sample Date/Time: 11/05/2015 09:12 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-7-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-7-A | | 240-206051 | 240-206089 | 11/10/2015 17:35 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-7-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-7-A | | 240-206256 | 240-205786 | 11/11/2015 13:44 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-7-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-7-A | | 240-206211 | 240-206227 | 11/11/2015 14:50 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-7-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-7-A | | 240-206105 | 240-205788 | 11/10/2015 23:46 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-7-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-7-B | | 240-206283 | 240-205809 | 11/11/2015 15:06 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-7 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-8** **Client ID: F-VB-UST-8-14**

Sample Date/Time: 11/05/2015 09:14 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-8-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-8-A | | 240-206051 | 240-206089 | 11/10/2015 17:56 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-D-8-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-D-8-A | | 240-206256 | 240-205786 | 11/11/2015 14:08 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-8-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-8-A | | 240-206211 | 240-206227 | 11/11/2015 15:27 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-E-8-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-E-8-A | | 240-206105 | 240-205788 | 11/11/2015 00:17 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-D-8-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-D-8-B | | 240-206283 | 240-205809 | 11/11/2015 15:23 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-8 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-9** **Client ID: F-VB-UST-9-14**

Sample Date/Time: 11/05/2015 09:16 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-9-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-9-A | | 240-206051 | 240-206089 | 11/10/2015 18:18 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-9-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-9-A | | 240-206256 | 240-205786 | 11/11/2015 14:33 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-9-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-9-A | | 240-206211 | 240-206227 | 11/11/2015 17:19 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-9-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-9-A | | 240-206105 | 240-205788 | 11/11/2015 00:48 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-9-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-9-B | | 240-206283 | 240-205809 | 11/11/2015 15:39 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-9 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-10** **Client ID: F-VB-UST-10-14**

Sample Date/Time: 11/05/2015 09:18 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-10-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-10-A | | 240-206051 | 240-206089 | 11/10/2015 18:39 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-10-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-10-A | | 240-206256 | 240-205786 | 11/11/2015 14:57 | 1 | TAL CAN | JMG |
| P:5030A | 240-57597-C-10-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-10-A | | 240-206211 | 240-206227 | 11/11/2015 17:57 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-10-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-10-A | | 240-206105 | 240-205788 | 11/11/2015 01:19 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-10-B | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-10-B | | 240-206283 | 240-205809 | 11/11/2015 15:55 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-10 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-10 DU** **Client ID: F-VB-UST-10-14**

Sample Date/Time: 11/05/2015 09:18 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| A:Moisture | 240-57597-A-10 DU | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-11** **Client ID: F-VB-UST-11-14**

Sample Date/Time: 11/05/2015 09:20 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-11-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-11-A | | 240-206051 | 240-206089 | 11/10/2015 19:01 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-11-B | | 240-206233 | 240-205866 | 11/09/2015 12:23 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-11-B | | 240-206233 | 240-205866 | 11/11/2015 14:50 | 1 | TAL CAN | MRU |
| P:5030A | 240-57597-C-11-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-11-A | | 240-206211 | 240-206227 | 11/11/2015 18:34 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-E-11-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-E-11-A | | 240-206105 | 240-205788 | 11/11/2015 01:50 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-D-11-A | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-D-11-A | | 240-206283 | 240-205809 | 11/11/2015 16:12 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-11 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-12** **Client ID: F-VB-UST-12-14**

Sample Date/Time: 11/05/2015 09:22 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-12-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-12-A | | 240-206051 | 240-206089 | 11/10/2015 19:22 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-12-B | | 240-206233 | 240-205866 | 11/09/2015 12:23 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-12-B | | 240-206233 | 240-205866 | 11/11/2015 15:14 | 1 | TAL CAN | MRU |
| P:5030A | 240-57597-C-12-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-12-A | | 240-206211 | 240-206227 | 11/11/2015 19:12 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-E-12-A | | 240-206356 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-E-12-A | | 240-206356 | 240-205788 | 11/11/2015 19:03 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-D-12-A | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-D-12-A | | 240-206283 | 240-205809 | 11/11/2015 16:28 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-12 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-13** **Client ID: F-VB-UST-13-14**

Sample Date/Time: 11/05/2015 09:24 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-13-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-13-A | | 240-206051 | 240-206089 | 11/10/2015 19:44 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-E-13-B | | 240-206233 | 240-205866 | 11/09/2015 12:23 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-E-13-B | | 240-206233 | 240-205866 | 11/11/2015 15:39 | 1 | TAL CAN | MRU |
| P:5030A | 240-57597-C-13-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-13-A | | 240-206211 | 240-206227 | 11/11/2015 19:48 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-E-13-A | | 240-206356 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-E-13-A | | 240-206356 | 240-205788 | 11/11/2015 19:34 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-D-13-A | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-D-13-A | | 240-206283 | 240-205809 | 11/11/2015 16:44 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-13 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-14** **Client ID: F-VB-UST-14-14**

Sample Date/Time: 11/05/2015 09:26 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-14-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-14-A | | 240-206051 | 240-206089 | 11/10/2015 20:05 | 1 | TAL CAN | TJL2 |
| P:3540C | 240-57597-D-14-B | | 240-206233 | 240-205866 | 11/09/2015 12:23 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-D-14-B | | 240-206233 | 240-205866 | 11/11/2015 16:03 | 1 | TAL CAN | MRU |
| P:5030A | 240-57597-C-14-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-14-A | | 240-206211 | 240-206227 | 11/11/2015 20:25 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-14-A | | 240-206356 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-14-A | | 240-206356 | 240-205788 | 11/11/2015 20:06 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-14-A | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-14-A | | 240-206283 | 240-205809 | 11/11/2015 17:01 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-14 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

**Lab ID: 240-57597-15** **Client ID: F-VB-UST-DUP**

Sample Date/Time: 11/05/2015 00:00 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-15-B | | 240-206267 | 240-206305 | 11/11/2015 11:03 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-15-B | | 240-206267 | 240-206305 | 11/11/2015 13:09 | 1 | TAL CAN | SAM |
| P:3540C | 240-57597-D-15-D | | 240-206233 | 240-205866 | 11/09/2015 12:23 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-D-15-D | | 240-206233 | 240-205866 | 11/11/2015 16:28 | 1 | TAL CAN | MRU |
| P:5030A | 240-57597-C-15-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-15-A | | 240-206211 | 240-206227 | 11/11/2015 21:03 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-15-A | | 240-206356 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-15-A | | 240-206356 | 240-205788 | 11/11/2015 20:37 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-15-A | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-15-A | | 240-206283 | 240-205809 | 11/11/2015 18:06 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57597-A-15 | | 240-205828 | | 11/09/2015 10:19 | 1 | TAL CAN | LCN |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-15 MS** **Client ID: F-VB-UST-DUP**

Sample Date/Time: 11/05/2015 00:00 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-15-C MS | | 240-206267 | 240-206305 | 11/11/2015 11:03 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-15-C MS | | 240-206267 | 240-206305 | 11/11/2015 13:31 | 1 | TAL CAN | SAM |
| P:3540C | 240-57597-D-15-E MS | | 240-206233 | 240-205866 | 11/09/2015 12:23 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-D-15-E MS | | 240-206233 | 240-205866 | 11/11/2015 16:52 | 1 | TAL CAN | MRU |
| P:5030A | 240-57597-C-15-B MS | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-15-B MS | | 240-206211 | 240-206227 | 11/11/2015 21:40 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-15-B MS | | 240-206356 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-15-B MS | | 240-206356 | 240-205788 | 11/11/2015 21:08 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-15-B MS | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-15-B MS | | 240-206283 | 240-205809 | 11/11/2015 18:22 | 1 | TAL CAN | LSH |

**Lab ID: 240-57597-15 MSD** **Client ID: F-VB-UST-DUP**

Sample Date/Time: 11/05/2015 00:00 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-15-D MSD | | 240-206267 | 240-206305 | 11/11/2015 11:03 | 1 | TAL CAN | SAM |
| A:8260C | 240-57597-B-15-D MSD | | 240-206267 | 240-206305 | 11/11/2015 13:52 | 1 | TAL CAN | SAM |
| P:3540C | 240-57597-D-15-F MSD | | 240-206233 | 240-205866 | 11/09/2015 12:23 | 1 | TAL CAN | YDN |
| A:8270D | 240-57597-D-15-F MSD | | 240-206233 | 240-205866 | 11/11/2015 17:17 | 1 | TAL CAN | MRU |
| P:5030A | 240-57597-C-15-C MSD | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | 240-57597-C-15-C MSD | | 240-206211 | 240-206227 | 11/11/2015 22:17 | 1 | TAL CAN | RTR |
| P:3540C | 240-57597-D-15-C MSD | | 240-206356 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | 240-57597-D-15-C MSD | | 240-206356 | 240-205788 | 11/11/2015 21:39 | 1 | TAL CAN | DEB |
| P:3540C | 240-57597-E-15-C MSD | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | 240-57597-E-15-C MSD | | 240-206283 | 240-205809 | 11/11/2015 18:39 | 1 | TAL CAN | LSH |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: 240-57597-16** **Client ID: TB-110515**

Sample Date/Time: 11/05/2015 00:00 Received Date/Time: 11/06/2015 10:20

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57597-B-16 | | 240-206461 | | 11/12/2015 12:33 | 1 | TAL CAN | LEE |
| A:8260C | 240-57597-B-16 | | 240-206461 | | 11/12/2015 12:33 | 1 | TAL CAN | LEE |

**Lab ID: MB** **Client ID: N/A**

Sample Date/Time: N/A Received Date/Time: N/A

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | MB 240-206089/1-A | | 240-206051 | 240-206089 | 11/10/2015 11:24 | 1 | TAL CAN | SAM |
| A:8260C | MB 240-206089/1-A | | 240-206051 | 240-206089 | 11/10/2015 12:56 | 1 | TAL CAN | TJL2 |
| A:8260C | MB 240-206267/6 | | 240-206267 | | 11/11/2015 11:43 | 1 | TAL CAN | SAM |
| P:5030C | MB 240-206461/6 | | 240-206461 | | 11/12/2015 12:11 | 1 | TAL CAN | LEE |
| A:8260C | MB 240-206461/6 | | 240-206461 | | 11/12/2015 12:11 | 1 | TAL CAN | LEE |
| P:3540C | MB 240-205866/8-A | | 240-206233 | 240-205866 | 11/09/2015 12:23 | 1 | TAL CAN | YDN |
| A:8270D | MB 240-205866/8-A | | 240-206233 | 240-205866 | 11/11/2015 09:54 | 1 | TAL CAN | MRU |
| P:3540C | MB 240-205786/23-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | MB 240-205786/23-A | | 240-206256 | 240-205786 | 11/11/2015 10:27 | 1 | TAL CAN | JMG |
| P:5030A | MB 240-206227/1-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | MB 240-206227/1-A | | 240-206211 | 240-206227 | 11/11/2015 09:56 | 1 | TAL CAN | RTR |
| P:5030A | MB 240-206444/1-A | | 240-206404 | 240-206444 | 11/12/2015 05:59 | 1 | TAL CAN | RTR |
| A:8015C | MB 240-206444/1-A | | 240-206404 | 240-206444 | 11/12/2015 08:03 | 1 | TAL CAN | RTR |
| P:3540C | MB 240-205788/22-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | MB 240-205788/22-A | | 240-206105 | 240-205788 | 11/10/2015 16:59 | 1 | TAL CAN | DEB |
| P:3540C | MB 240-205809/21-A | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | MB 240-205809/21-A | | 240-206283 | 240-205809 | 11/11/2015 17:17 | 1 | TAL CAN | LSH |

Client: Tetra Tech, Inc. Job Number: 240-57597-1

## Laboratory Chronicle

**Lab ID: LCS** **Client ID: N/A**

Sample Date/Time: N/A Received Date/Time: N/A

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| A:8260C | LCS 240-206051/5 | | 240-206051 | | 11/10/2015 11:30 | 1 | TAL CAN | TJL2 |
| A:8260C | LCS 240-206267/5 | | 240-206267 | | 11/11/2015 11:22 | 1 | TAL CAN | SAM |
| P:5030C | LCS 240-206461/4 | | 240-206461 | | 11/12/2015 11:03 | 1 | TAL CAN | LEE |
| A:8260C | LCS 240-206461/4 | | 240-206461 | | 11/12/2015 11:03 | 1 | TAL CAN | LEE |
| P:3540C | LCS 240-205866/9-A | | 240-206233 | 240-205866 | 11/09/2015 12:23 | 1 | TAL CAN | YDN |
| A:8270D | LCS 240-205866/9-A | | 240-206233 | 240-205866 | 11/11/2015 10:19 | 1 | TAL CAN | MRU |
| P:3540C | LCS 240-205786/24-A | | 240-206256 | 240-205786 | 11/09/2015 06:12 | 1 | TAL CAN | YDN |
| A:8270D | LCS 240-205786/24-A | | 240-206256 | 240-205786 | 11/11/2015 10:51 | 1 | TAL CAN | JMG |
| P:5030A | LCS 240-206227/2-A | | 240-206211 | 240-206227 | 11/11/2015 06:58 | 1 | TAL CAN | RTR |
| A:8015C | LCS 240-206227/2-A | | 240-206211 | 240-206227 | 11/11/2015 10:31 | 1 | TAL CAN | RTR |
| P:5030A | LCS 240-206444/2-A | | 240-206404 | 240-206444 | 11/12/2015 05:59 | 1 | TAL CAN | RTR |
| A:8015C | LCS 240-206444/2-A | | 240-206404 | 240-206444 | 11/12/2015 08:40 | 1 | TAL CAN | RTR |
| P:3540C | LCS 240-205788/23-A | | 240-206105 | 240-205788 | 11/09/2015 09:40 | 1 | TAL CAN | YDN |
| A:8015C | LCS 240-205788/23-A | | 240-206105 | 240-205788 | 11/10/2015 17:31 | 1 | TAL CAN | DEB |
| P:3540C | LCS 240-205809/22-A | | 240-206283 | 240-205809 | 11/09/2015 09:53 | 1 | TAL CAN | YDN |
| A:8082A | LCS 240-205809/22-A | | 240-206283 | 240-205809 | 11/11/2015 17:33 | 1 | TAL CAN | LSH |

**Lab ID: MS** **Client ID: N/A**

Sample Date/Time: 11/05/2015 12:10 Received Date/Time: 11/06/2015 09:00

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57564-B-1 MS | | 240-206461 | | 11/12/2015 20:05 | 250 | TAL CAN | LEE |
| A:8260C | 240-57564-B-1 MS | | 240-206461 | | 11/12/2015 20:05 | 250 | TAL CAN | LEE |

**Lab ID: MSD** **Client ID: N/A**

Sample Date/Time: 11/05/2015 12:10 Received Date/Time: 11/06/2015 09:00

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57564-B-1 MSD | | 240-206461 | | 11/12/2015 20:27 | 250 | TAL CAN | LEE |
| A:8260C | 240-57564-B-1 MSD | | 240-206461 | | 11/12/2015 20:27 | 250 | TAL CAN | LEE |

**Lab References:**
TAL CAN = TestAmerica Canton

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **ex10PPMSPK_00018** | 04/30/19 | 08/21/15 | MEOH, Lot 147358 | 1000 mL | ex1016/1260st_00001 | 3.8 mL | Aroclor-1016 | 10.0012 ug/mL |
| | | | | | | | Aroclor-1260 | 10.0062 ug/mL |
| | | | | | ex1016/1260st_00002 | 6.2 mL | Aroclor-1016 | 10.0012 ug/mL |
| | | | | | | | Aroclor-1260 | 10.0062 ug/mL |
| .ex1016/1260st_00001 | 04/30/19 | | Restek, Lot A092844 | | (Purchased Reagent) | | Aroclor-1016 | 1000 ug/mL |
| | | | | | | | Aroclor-1260 | 1000 ug/mL |
| .ex1016/1260st_00002 | 05/31/21 | | Restek, Lot A0109056 | | (Purchased Reagent) | | Aroclor-1016 | 1000.2 ug/mL |
| | | | | | | | Aroclor-1260 | 1001 ug/mL |
| **ex2/.2SURRW_00062** | 07/13/17 | 10/26/15 | MEOH, Lot 0000118655 | 2000 mL | EXPESTSURR_00001 | 2 mL | DCB Decachlorobiphenyl | 0.2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.2 ug/mL |
| .EXPESTSURR_00001 | 05/31/21 | | Restek, Lot A0108779 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **exBENZALDEHYD_00031** | 08/31/16 | 10/27/15 | MEOH, Lot 0000099494 | 200 mL | EXLIST1_S11_00001 | 4 mL | Atrazine | 40 ug/mL |
| | | | | | | | Benzaldehyde | 40 ug/mL |
| | | | | | | | Caprolactam | 40 ug/mL |
| .EXLIST1_S11_00001 | 08/31/16 | | Restek, Lot A0108989 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| **exBNASPIKE_00053** | 08/31/16 | 11/02/15 | MEOH, Lot 0000099494 | 200 mL | exLIST1_S1_00004 | 4 mL | 1,1'-Biphenyl | 20 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 20 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 20 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dioxane | 20 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 20 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 40 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 20 ug/mL |
| | | | | | | | 2-Chlorophenol | 20 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2-Methylphenol | 20 ug/mL |
| | | | | | | | 2-Nitroaniline | 20 ug/mL |
| | | | | | | | 2-Nitrophenol | 20 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 20 ug/mL |
| | | | | | | | 3-Methylphenol | 10 ug/mL |
| | | | | | | | 3-Nitroaniline | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 40 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 20 ug/mL |
| | | | | | | | 4-Chloroaniline | 20 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Methylphenol | 10 ug/mL |
| | | | | | | | 4-Nitroaniline | 20 ug/mL |
| | | | | | | | 4-Nitrophenol | 40 ug/mL |
| | | | | | | | Acenaphthene | 20 ug/mL |
| | | | | | | | Acenaphthylene | 20 ug/mL |
| | | | | | | | Acetophenone | 20 ug/mL |
| | | | | | | | Aniline | 20 ug/mL |
| | | | | | | | Anthracene | 20 ug/mL |
| | | | | | | | Azobenzene | 20 ug/mL |
| | | | | | | | Benzo[a]anthracene | 20 ug/mL |
| | | | | | | | Benzo[a]pyrene | 20 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 20 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 20 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 20 ug/mL |
| | | | | | | | Benzyl alcohol | 20 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 20 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 20 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 20 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 20 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 20 ug/mL |
| | | | | | | | Carbazole | 20 ug/mL |
| | | | | | | | Chrysene | 20 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 20 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 20 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 20 ug/mL |
| | | | | | | | Dibenzofuran | 20 ug/mL |
| | | | | | | | Diethyl phthalate | 20 ug/mL |
| | | | | | | | Dimethyl phthalate | 20 ug/mL |
| | | | | | | | Fluoranthene | 20 ug/mL |
| | | | | | | | Fluorene | 20 ug/mL |
| | | | | | | | Hexachlorobenzene | 20 ug/mL |
| | | | | | | | Hexachlorobutadiene | 20 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 20 ug/mL |
| | | | | | | | Hexachloroethane | 20 ug/mL |
| | | | | | | | Hexadecane | 20 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 20 ug/mL |
| | | | | | | | Isophorone | 20 ug/mL |
| | | | | | | | n-Decane | 20 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 40 ug/mL |
| | | | | | | | n-Octadecane | 20 ug/mL |
| | | | | | | | Naphthalene | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Nitrobenzene | 20 ug/mL |
| | | | | | | | Pentachlorophenol | 40 ug/mL |
| | | | | | | | Phenanthrene | 20 ug/mL |
| | | | | | | | Phenol | 20 ug/mL |
| | | | | | | | Pyrene | 20 ug/mL |
| | | | | | | | Pyridine | 20 ug/mL |
| | | | | | EXLIST1_S10_00001 | 4 mL | Benzoic acid | 40 ug/mL |
| | | | | | | | Indene | 40 ug/mL |
| | | | | | exLIST1_S9_00003 | 4 mL | 3,3'-Dichlorobenzidine | 40 ug/mL |
| | | | | | | | Benzidine | 40 ug/mL |
| .exLIST1_S1_00004 | 12/31/16 | | Restek, Lot A0111934 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Methylphenol | 500 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Methylphenol | 500 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| .EXLIST1_S10_00001 | 08/31/16 | | Restek, Lot A0108988 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| .exLIST1_S9_00003 | 01/31/17 | | Restek, Lot A0112567 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **exBNASURR W_00028** | 02/22/18 | 10/15/15 | MEOH, Lot 0000099494 | 2000 mL | SMLIST1 SURR_00003 | 8 mL | 2,4,6-Tribromophenol (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 20 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 20 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 20 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 20 ug/mL |
| .SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **exDIESELspkw_00068** | 08/27/17 | 10/05/15 | ACETONE, Lot 0000090987 | 200 mL | exDIESELstd_00003 | 10 mL | Diesel | 2521.45 ug/mL |
| .exDIESELstd_00003 | 04/30/22 | | Restek, Lot A0109732 | | (Purchased Reagent) | | Diesel | 50429 ug/mL |
| **exOTPSURRW_00006** | 12/31/18 | 10/17/15 | MEOH, Lot 0000099494 | 1000 mL | exOTPstd_00002 | 20 mL | o-Terphenyl (Surr) | 200 ug/mL |
| .exOTPstd_00002 | 12/31/18 | | Restek, Lot A0110923 | | (Purchased Reagent) | | o-Terphenyl (Surr) | 10000 ug/mL |
| **Gas Surr Cali_00014** | 11/02/15 | 08/07/15 | MeoH, Lot 1823933 | 10 mL | VGGASURROGATE_00032 | 1 mL | Trifluorotoluene (Surr) | 5 ug/mL |
| .VGGASURROGATE_00032 | 08/22/15 | 05/22/15 | MEOH, Lot 49909 | 50 mL | VGSURRSTKGAS_00019 | 1.25 mL | Trifluorotoluene (Surr) | 50 ug/mL |
| ..VGSURRSTKGAS_00019 | 08/31/16 | | ULTRA SCIENTIFIC, Lot CE-1903A | | (Purchased Reagent) | | Trifluorotoluene (Surr) | 2000 ug/mL |
| **SG1248@.05ppm_00018** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 0.5 mL | PCB-1248 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.05 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.1PPM_00015** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 1 mL | PCB-1248 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.1 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.2ppm_00015** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 2 mL | PCB-1248 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.5ppm_00020** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 5 mL | PCB-1248 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.5 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.5ppm_00020** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 5 mL | Aroclor-1248 | 0.5 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | Aroclor-1248 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | Aroclor-1248 | 1000 ug/mL |
| **SG1248@1.0ppm_00014** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 10 mL | PCB-1248 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@2.0ppm_00014** | 04/22/16 | 08/24/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 20 mL | PCB-1248 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1660@.05PPM_00015** | 03/02/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1016/1260_00017 | 0.5 mL | PCB-1016 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.05 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.125 mL | DCB Decachlorobiphenyl | 0.0025 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.0025 ug/mL |
| .SG1016/1260_00017 | 03/02/16 | 09/02/15 | HEXANE, Lot 000017549 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@0.2ppm_00016** | 04/22/16 | 11/13/15 | HEXANE, Lot 154006 | 100 mL | SG1660ICV@10_00008 | 2 mL | PCB-1016 Peak 1 | 0.00002 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.00002 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.00002 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.00002 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.00002 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.00002 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.00002 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.00002 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.00002 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.00002 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.5 mL | DCB Decachlorobiphenyl | 0.01 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.01 ug/mL |
| .SG1660ICV@10_00008 | 05/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 10 mL | SG 1016 ICV_00007 | 1 mL | PCB-1016 Peak 1 | 0.001 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.001 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.001 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.001 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.001 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | SG 1260 ICV_00008 | 1 mL | PCB-1260 Peak 1 | 0.001 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.001 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.001 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.001 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.001 ug/mL |
| ..SG 1016 ICV_00007 | 02/28/20 | | ultra, Lot CJ-0176 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 10 ug/L |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/L |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/L |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/L |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/L |
| ..SG 1260 ICV_00008 | 02/28/17 | | Ultra, Lot CG-3829A | | (Purchased Reagent) | | PCB-1260 Peak 1 | 10 ug/L |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/L |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/L |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/L |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/L |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@0.5PPM_00056** | 03/02/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1016/1260_00017 | 5 mL | PCB-1016 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.5 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 1.25 mL | DCB Decachlorobiphenyl | 0.025 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.025 ug/mL |
| .SG1016/1260_00017 | 03/02/16 | 09/02/15 | HEXANE, Lot 000017549 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@0.5PPM_00056** | 03/02/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1016/1260_00017 | 5 mL | Aroclor-1016 | 0.5 ug/mL |
| | | | | | | | Aroclor-1260 | 0.5 ug/mL |
| .SG1016/1260_00017 | 03/02/16 | 09/02/15 | HEXANE, Lot 000017549 | 100 mL | SG1016/1260MX_00009 | 1 mL | Aroclor-1016 | 10 ug/mL |
| | | | | | | | Aroclor-1260 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | Aroclor-1016 | 1000 ug/mL |
| | | | | | | | Aroclor-1260 | 1000 ug/mL |
| **SG1660@1.0PPM_00013** | 03/02/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1016/1260_00017 | 10 mL | PCB-1016 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 2.5 mL | DCB Decachlorobiphenyl | 0.05 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.05 ug/mL |
| .SG1016/1260_00017 | 03/02/16 | 09/02/15 | HEXANE, Lot 000017549 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@2.0PPM_00014** | 04/22/16 | 11/13/15 | HEXANE, Lot 00000101243 | 100 mL | SG1660ICV@10_00008 | 20 mL | PCB-1016 Peak 1 | 0.0002 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.0002 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.0002 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.0002 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.0002 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.0002 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.0002 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.0002 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.0002 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.0002 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 5 mL | DCB Decachlorobiphenyl | 0.1 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.1 ug/mL |
| .SG1660ICV@10_00008 | 05/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 10 mL | SG 1016 ICV_00007 | 1 mL | PCB-1016 Peak 1 | 0.001 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.001 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.001 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.001 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.001 ug/mL |
| | | | | | SG 1260 ICV_00008 | 1 mL | PCB-1260 Peak 1 | 0.001 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.001 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.001 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.001 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.001 ug/mL |
| ..SG 1016 ICV_00007 | 02/28/20 | | ultra, Lot CJ-0176 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 10 ug/L |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/L |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/L |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/L |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/L |
| ..SG 1260 ICV_00008 | 02/28/17 | | Ultra, Lot CG-3829A | | (Purchased Reagent) | | PCB-1260 Peak 1 | 10 ug/L |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/L |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/L |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/L |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/L |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660ICV@1.0_00014** | 04/22/16 | 10/28/15 | HEXANE, Lot 0000101243 | 10 mL | SGPCBIS STOCk_00003 | 0.05 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1660ICV@1.0_00014** | 04/22/16 | 10/28/15 | HEXANE, Lot 0000101243 | 10 mL | SG1660ICV@100_00005 | 0.1 mL | Aroclor-1016 | 1 ug/mL |
| | | | | | | | Aroclor-1260 | 1 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.25 mL | DCB Decachlorobiphenyl | 0.05 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.05 ug/mL |
| .SG1660ICV@100_00005 | 10/28/16 | 10/28/15 | HEXANE, Lot 0000101243 | 10 mL | SG 1660 ICV_00012 | 1 mL | Aroclor-1016 | 100 ug/mL |
| | | | | | | | Aroclor-1260 | 100 ug/mL |
| ..SG 1660 ICV_00012 | 01/30/22 | | restek, Lot a0114674 | | (Purchased Reagent) | | Aroclor-1016 | 1000 ug/mL |
| | | | | | | | Aroclor-1260 | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660STD@0.1_00015** | 03/02/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1016/1260_00017 | 1 mL | PCB-1016 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.1 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.25 mL | DCB Decachlorobiphenyl | 0.005 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.005 ug/mL |
| .SG1016/1260_00017 | 03/02/16 | 09/02/15 | HEXANE, Lot 000017549 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG2154@0.05PP_00010** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 0.5 mL | PCB-1221 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.05 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@0.2PPM_00012** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 2 mL | PCB-1221 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.2 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@0.5PPM_00021** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 5 mL | PCB-1221 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.5 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@0.5PPM_00021** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 5 mL | Aroclor-1221 | 0.5 ug/mL |
| | | | | | | | Aroclor-1254 | 0.5 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | Aroclor-1221 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Aroclor-1254 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | Aroclor-1221 | 1000 ug/mL |
| | | | | | | | Aroclor-1254 | 1000 ug/mL |
| **SG2154@1.0PPM_00014** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 10 mL | PCB-1221 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@2.0PPM_00011** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 20 mL | PCB-1221 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 2 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@0.1PPM_00013** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 1 mL | PCB-1221 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.1 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@.05PPM_00004** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00003 | 0.5 mL | PCB-1232 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 50 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00003 | 09/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/62 STK_00001 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SG32/62 STK_00001 | 04/30/21 | | Restelk, Lot A0108264 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@.2PPM_00005** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 2 mL | PCB-1232 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 200 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@0.1PPM_00006** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 1 mL | PCB-1232 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 100 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@0.5PPM_00010** | 04/22/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00003 | 5 mL | PCB-1232 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 500 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1262 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 500 ug/L |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00003 | 09/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/62 STK_00001 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SG32/62 STK_00001 | 04/30/21 | | Restelk, Lot A0108264 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@0.5PPM_00010** | 04/22/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00003 | 5 mL | Aroclor-1232 | 500 ug/L |
| | | | | | | | PCB-1262 | 500 ug/L |
| .SG32/6210PPM_00003 | 09/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/62 STK_00001 | 1 mL | Aroclor-1232 | 10000 ug/L |
| | | | | | | | PCB-1262 | 10000 ug/L |
| ..SG32/62 STK_00001 | 04/30/21 | | Restelk, Lot A0108264 | | (Purchased Reagent) | | Aroclor-1232 | 1000 ug/mL |
| | | | | | | | PCB-1262 | 1000 ug/mL |
| **SG3262@1.0PPM_00005** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 10 mL | PCB-1232 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@2.0ppm_00004** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 20 mL | PCB-1232 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 2000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG42/68@1.0PP_00008** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 10 mL | PCB-1242 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@.05PPM_00005** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 0.5 mL | PCB-1242 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 50 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1268 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 50 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@.1PPM_00006** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 1 mL | PCB-1242 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 100 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@.2PPM_00005** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 2 mL | PCB-1242 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 200 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@0.5PPM_00007** | 12/24/15 | 06/24/15 | HEXANE, Lot 0000086560 | 100 mL | SG42/6810PPM_00002 | 5 mL | PCB-1242 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 500 ug/L |

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1242 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 500 ug/L |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@0.5PPM_00011** | 04/07/16 | 10/07/15 | HEXANE, Lot 154006 | 100 mL | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@0.5PPM_00011** | 04/07/16 | 10/07/15 | HEXANE, Lot 154006 | 100 mL | SG42/6810PPM_00003 | 5 mL | Aroclor-1242 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 500 ug/L |
| | | | | | | | PCB-1268 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 500 ug/L |
| .SG42/6810PPM_00003 | 10/07/16 | 10/07/15 | HEXANE, Lot 154006 | 100 mL | SG 42/68 STK_00001 | 1 mL | Aroclor-1242 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 1 | 10000 ug/L |

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | | | PCB-1268 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG 42/68 STK_00001 | 04/30/21 | | RESTEK, Lot A0108269 | | (Purchased Reagent) | | Aroclor-1242 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1268 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| **SG4268@2.0PPM_00004** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 20 mL | PCB-1242 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 2000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SGDROCAL_00020** | 10/05/16 | 10/05/15 | MECL2, Lot EXDCM_00187 | 25 mL | SGDROCALSTK_00019 | 5 mL | Diesel | 10002.5 ug/mL |
| | | | | | SGDROSURROTP_00025 | 1 mL | o-Terphenyl (Surr) | 800 ug/mL |
| | | | | | SGDROSURROTP_00026 | 1 mL | o-Terphenyl (Surr) | 800 ug/mL |
| .SGDROCALSTK_00019 | 04/30/22 | | Restek, Lot A0109723 | | (Purchased Reagent) | | Diesel | 50005.8 ug/mL |
| .SGDROSURROTP_00025 | 09/30/18 | | Restek, Lot A0109488 | | (Purchased Reagent) | | o-Terphenyl (Surr) | 10062.7 ug/mL |
| .SGDROSURROTP_00026 | 09/30/18 | | Restek, Lot A0109488 | | (Purchased Reagent) | | o-Terphenyl (Surr) | 10062.7 ug/mL |
| **SGDROICV_00015** | 08/26/16 | 08/26/15 | MECL2, Lot 2091477 | 25 mL | SGDROICVSTK_00008 | 1 mL | Diesel | 2006 ug/mL |
| .SGDROICVSTK_00008 | 08/31/18 | | Ultra, Lot CL-2929 | | (Purchased Reagent) | | Diesel | 50150 ug/mL |
| **SGPCBISTD_00003** | 04/22/16 | 10/13/15 | HEXANE, Lot 154006 | 100 mL | SGPCBIS STOCk_00003 | 10 mL | 1-Bromo-2-nitrobenzene | 1 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SMIS80PPMW_00013** | 07/10/16 | 07/10/15 | MECL2, Lot 0000103549 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| .SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| **SMLIST1 L1 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1PAH ST_00004 | 2 uL | 1,1'-Biphenyl | 0.1 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 0.1 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 0.1 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 0.1 ug/mL |
| | | | | | | | Acenaphthene | 0.1 ug/mL |
| | | | | | | | Acenaphthylene | 0.1 ug/mL |
| | | | | | | | Anthracene | 0.1 ug/mL |
| | | | | | | | Benzo[a]anthracene | 0.1 ug/mL |
| | | | | | | | Benzo[a]pyrene | 0.1 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[b]fluoranthene | 0.1 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 0.1 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 0.1 ug/mL |
| | | | | | | | Chrysene | 0.1 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 0.1 ug/mL |
| | | | | | | | Dibenzofuran | 0.1 ug/mL |
| | | | | | | | Fluoranthene | 0.1 ug/mL |
| | | | | | | | Fluorene | 0.1 ug/mL |
| | | | | | | | Hexachlorobenzene | 0.1 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 0.1 ug/mL |
| | | | | | | | Naphthalene | 0.1 ug/mL |
| | | | | | | | Phenanthrene | 0.1 ug/mL |
| | | | | | | | Pyrene | 0.1 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 0.1 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 0.1 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 0.1 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1PAH ST_00004 | 05/31/16 | 01/26/15 | MECL2, Lot 0000093871 | 3 mL | SMLIST1 PAH_00004 | 300 uL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 60 uL | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 PAH_00004 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L2 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 10 uL | 1,1'-Biphenyl | 0.5 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 0.5 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 0.5 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,4-Dioxane | 0.5 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 0.5 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 0.5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,4,5-Trichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 0.5 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 1 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 0.5 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 0.5 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 0.5 ug/mL |
| | | | | | | | 2-Chlorophenol | 0.5 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 0.5 ug/mL |
| | | | | | | | 2-Methylphenol | 0.5 ug/mL |
| | | | | | | | 2-Nitroaniline | 0.5 ug/mL |
| | | | | | | | 2-Nitrophenol | 0.5 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 0.5 ug/mL |
| | | | | | | | 3-Nitroaniline | 0.5 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 1 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 0.5 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 0.5 ug/mL |
| | | | | | | | 4-Chloroaniline | 0.5 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 0.5 ug/mL |
| | | | | | | | 4-Nitroaniline | 0.5 ug/mL |
| | | | | | | | 4-Nitrophenol | 1 ug/mL |
| | | | | | | | Acenaphthene | 0.5 ug/mL |
| | | | | | | | Acenaphthylene | 0.5 ug/mL |
| | | | | | | | Acetophenone | 0.5 ug/mL |
| | | | | | | | Aniline | 0.5 ug/mL |
| | | | | | | | Anthracene | 0.5 ug/mL |
| | | | | | | | Azobenzene | 0.5 ug/mL |
| | | | | | | | Benzo[a]anthracene | 0.5 ug/mL |
| | | | | | | | Benzo[a]pyrene | 0.5 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 0.5 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 0.5 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 0.5 ug/mL |
| | | | | | | | Benzyl alcohol | 0.5 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 0.5 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 0.5 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 0.5 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 0.5 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 0.5 ug/mL |
| | | | | | | | Carbazole | 0.5 ug/mL |
| | | | | | | | Chrysene | 0.5 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 0.5 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 0.5 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 0.5 ug/mL |
| | | | | | | | Dibenzofuran | 0.5 ug/mL |
| | | | | | | | Diethyl phthalate | 0.5 ug/mL |
| | | | | | | | Dimethyl phthalate | 0.5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Fluoranthene | 0.5 ug/mL |
| | | | | | | | Fluorene | 0.5 ug/mL |
| | | | | | | | Hexachlorobenzene | 0.5 ug/mL |
| | | | | | | | Hexachlorobutadiene | 0.5 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 0.5 ug/mL |
| | | | | | | | Hexachloroethane | 0.5 ug/mL |
| | | | | | | | Hexadecane | 0.5 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 0.5 ug/mL |
| | | | | | | | Isophorone | 0.5 ug/mL |
| | | | | | | | n-Decane | 0.5 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 0.5 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 0.5 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 1 ug/mL |
| | | | | | | | n-Octadecane | 0.5 ug/mL |
| | | | | | | | Naphthalene | 0.5 ug/mL |
| | | | | | | | Nitrobenzene | 0.5 ug/mL |
| | | | | | | | Pentachlorophenol | 1 ug/mL |
| | | | | | | | Phenanthrene | 0.5 ug/mL |
| | | | | | | | Phenol | 0.5 ug/mL |
| | | | | | | | Pyrene | 0.5 ug/mL |
| | | | | | | | Pyridine | 0.5 ug/mL |
| | | | | | | | Benzoic acid | 1 ug/mL |
| | | | | | | | Indene | 1 ug/mL |
| | | | | | | | Atrazine | 1 ug/mL |
| | | | | | | | Benzaldehyde | 1 ug/mL |
| | | | | | | | Caprolactam | 1 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 1 ug/mL |
| | | | | | | | Benzidine | 1 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 0.5 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 0.5 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 0.5 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 0.5 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 0.5 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 0.5 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L3 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 20 uL | 1,1'-Biphenyl | 1 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1 ug/mL |
| | | | | | | | 1,4-Dioxane | 1 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1 ug/mL |
| | | | | | | | 2-Chlorophenol | 1 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1 ug/mL |
| | | | | | | | 2-Methylphenol | 1 ug/mL |
| | | | | | | | 2-Nitroaniline | 1 ug/mL |
| | | | | | | | 2-Nitrophenol | 1 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1 ug/mL |
| | | | | | | | 3-Nitroaniline | 1 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1 ug/mL |
| | | | | | | | 4-Chloroaniline | 1 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1 ug/mL |
| | | | | | | | 4-Nitroaniline | 1 ug/mL |
| | | | | | | | 4-Nitrophenol | 2 ug/mL |
| | | | | | | | Acenaphthene | 1 ug/mL |
| | | | | | | | Acenaphthylene | 1 ug/mL |
| | | | | | | | Acetophenone | 1 ug/mL |
| | | | | | | | Aniline | 1 ug/mL |
| | | | | | | | Anthracene | 1 ug/mL |
| | | | | | | | Azobenzene | 1 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1 ug/mL |
| | | | | | | | Benzyl alcohol | 1 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1 ug/mL |
| | | | | | | | Carbazole | 1 ug/mL |
| | | | | | | | Chrysene | 1 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1 ug/mL |
| | | | | | | | Dibenzofuran | 1 ug/mL |
| | | | | | | | Diethyl phthalate | 1 ug/mL |
| | | | | | | | Dimethyl phthalate | 1 ug/mL |
| | | | | | | | Fluoranthene | 1 ug/mL |
| | | | | | | | Fluorene | 1 ug/mL |
| | | | | | | | Hexachlorobenzene | 1 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1 ug/mL |
| | | | | | | | Hexachloroethane | 1 ug/mL |
| | | | | | | | Hexadecane | 1 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1 ug/mL |
| | | | | | | | Isophorone | 1 ug/mL |
| | | | | | | | n-Decane | 1 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2 ug/mL |
| | | | | | | | n-Octadecane | 1 ug/mL |
| | | | | | | | Naphthalene | 1 ug/mL |
| | | | | | | | Nitrobenzene | 1 ug/mL |
| | | | | | | | Pentachlorophenol | 2 ug/mL |
| | | | | | | | Phenanthrene | 1 ug/mL |
| | | | | | | | Phenol | 1 ug/mL |
| | | | | | | | Pyrene | 1 ug/mL |
| | | | | | | | Pyridine | 1 ug/mL |
| | | | | | | | Benzoic acid | 2 ug/mL |
| | | | | | | | Indene | 2 ug/mL |
| | | | | | | | Atrazine | 2 ug/mL |
| | | | | | | | Benzaldehyde | 2 ug/mL |
| | | | | | | | Caprolactam | 2 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 2 ug/mL |
| | | | | | | | Benzidine | 2 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 1 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 1 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 1 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 1 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 1 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 1 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L4 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 40 uL | 1,1'-Biphenyl | 2 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 2 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 2 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 2 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2 ug/mL |
| | | | | | | | 1,4-Dioxane | 2 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 2 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 2 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 2 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 2 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 2 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 2 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 4 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 2 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 2 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 2 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 2 ug/mL |
| | | | | | | | 2-Chlorophenol | 2 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 2 ug/mL |
| | | | | | | | 2-Methylphenol | 2 ug/mL |
| | | | | | | | 2-Nitroaniline | 2 ug/mL |
| | | | | | | | 2-Nitrophenol | 2 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 2 ug/mL |
| | | | | | | | 3-Nitroaniline | 2 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 4 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 2 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 2 ug/mL |
| | | | | | | | 4-Chloroaniline | 2 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 2 ug/mL |
| | | | | | | | 4-Nitroaniline | 2 ug/mL |
| | | | | | | | 4-Nitrophenol | 4 ug/mL |
| | | | | | | | Acenaphthene | 2 ug/mL |
| | | | | | | | Acenaphthylene | 2 ug/mL |
| | | | | | | | Acetophenone | 2 ug/mL |
| | | | | | | | Aniline | 2 ug/mL |
| | | | | | | | Anthracene | 2 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Azobenzene | 2 ug/mL |
| | | | | | | | Benzo[a]anthracene | 2 ug/mL |
| | | | | | | | Benzo[a]pyrene | 2 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 2 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 2 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 2 ug/mL |
| | | | | | | | Benzyl alcohol | 2 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 2 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 2 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 2 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 2 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 2 ug/mL |
| | | | | | | | Carbazole | 2 ug/mL |
| | | | | | | | Chrysene | 2 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 2 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 2 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 2 ug/mL |
| | | | | | | | Dibenzofuran | 2 ug/mL |
| | | | | | | | Diethyl phthalate | 2 ug/mL |
| | | | | | | | Dimethyl phthalate | 2 ug/mL |
| | | | | | | | Fluoranthene | 2 ug/mL |
| | | | | | | | Fluorene | 2 ug/mL |
| | | | | | | | Hexachlorobenzene | 2 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 2 ug/mL |
| | | | | | | | Hexachloroethane | 2 ug/mL |
| | | | | | | | Hexadecane | 2 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 2 ug/mL |
| | | | | | | | Isophorone | 2 ug/mL |
| | | | | | | | n-Decane | 2 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 2 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 2 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 4 ug/mL |
| | | | | | | | n-Octadecane | 2 ug/mL |
| | | | | | | | Naphthalene | 2 ug/mL |
| | | | | | | | Nitrobenzene | 2 ug/mL |
| | | | | | | | Pentachlorophenol | 4 ug/mL |
| | | | | | | | Phenanthrene | 2 ug/mL |
| | | | | | | | Phenol | 2 ug/mL |
| | | | | | | | Pyrene | 2 ug/mL |
| | | | | | | | Pyridine | 2 ug/mL |
| | | | | | | | Benzoic acid | 4 ug/mL |
| | | | | | | | Indene | 4 ug/mL |
| | | | | | | | Atrazine | 4 ug/mL |
| | | | | | | | Benzaldehyde | 4 ug/mL |
| | | | | | | | Caprolactam | 4 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 4 ug/mL |
| | | | | | | | Benzidine | 4 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 2 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 2 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 2 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 2 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 2 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 2 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L5 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 100 uL | 1,1'-Biphenyl | 5 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 5 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 5 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 5 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 5 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 5 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 5 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 5 ug/mL |
| | | | | | | | 1,4-Dioxane | 5 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 5 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 5 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 5 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 5 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 5 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 5 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 10 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 5 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 5 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 5 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 5 ug/mL |
| | | | | | | | 2-Chlorophenol | 5 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 5 ug/mL |
| | | | | | | | 2-Methylphenol | 5 ug/mL |
| | | | | | | | 2-Nitroaniline | 5 ug/mL |
| | | | | | | | 2-Nitrophenol | 5 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 5 ug/mL |
| | | | | | | | 3-Nitroaniline | 5 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 10 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Chloro-3-methylphenol | 5 ug/mL |
| | | | | | | | 4-Chloroaniline | 5 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 5 ug/mL |
| | | | | | | | 4-Nitroaniline | 5 ug/mL |
| | | | | | | | 4-Nitrophenol | 10 ug/mL |
| | | | | | | | Acenaphthene | 5 ug/mL |
| | | | | | | | Acenaphthylene | 5 ug/mL |
| | | | | | | | Acetophenone | 5 ug/mL |
| | | | | | | | Aniline | 5 ug/mL |
| | | | | | | | Anthracene | 5 ug/mL |
| | | | | | | | Azobenzene | 5 ug/mL |
| | | | | | | | Benzo[a]anthracene | 5 ug/mL |
| | | | | | | | Benzo[a]pyrene | 5 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 5 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 5 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 5 ug/mL |
| | | | | | | | Benzyl alcohol | 5 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 5 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 5 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 5 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 5 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 5 ug/mL |
| | | | | | | | Carbazole | 5 ug/mL |
| | | | | | | | Chrysene | 5 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 5 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 5 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 5 ug/mL |
| | | | | | | | Dibenzofuran | 5 ug/mL |
| | | | | | | | Diethyl phthalate | 5 ug/mL |
| | | | | | | | Dimethyl phthalate | 5 ug/mL |
| | | | | | | | Fluoranthene | 5 ug/mL |
| | | | | | | | Fluorene | 5 ug/mL |
| | | | | | | | Hexachlorobenzene | 5 ug/mL |
| | | | | | | | Hexachlorobutadiene | 5 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 5 ug/mL |
| | | | | | | | Hexachloroethane | 5 ug/mL |
| | | | | | | | Hexadecane | 5 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 5 ug/mL |
| | | | | | | | Isophorone | 5 ug/mL |
| | | | | | | | n-Decane | 5 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 5 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 5 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 10 ug/mL |
| | | | | | | | n-Octadecane | 5 ug/mL |
| | | | | | | | Naphthalene | 5 ug/mL |
| | | | | | | | Nitrobenzene | 5 ug/mL |
| | | | | | | | Pentachlorophenol | 10 ug/mL |
| | | | | | | | Phenanthrene | 5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Phenol | 5 ug/mL |
| | | | | | | | Pyrene | 5 ug/mL |
| | | | | | | | Pyridine | 5 ug/mL |
| | | | | | | | Benzoic acid | 10 ug/mL |
| | | | | | | | Indene | 10 ug/mL |
| | | | | | | | Atrazine | 10 ug/mL |
| | | | | | | | Benzaldehyde | 10 ug/mL |
| | | | | | | | Caprolactam | 10 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 10 ug/mL |
| | | | | | | | Benzidine | 10 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 5 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L6 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 200 uL | 1,1'-Biphenyl | 10 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 10 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 10 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 10 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 10 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 10 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 10 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 10 ug/mL |
| | | | | | | | 1,4-Dioxane | 10 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 10 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 10 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 10 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 10 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Chloronaphthalene | 10 ug/mL |
| | | | | | | | 2-Chlorophenol | 10 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 10 ug/mL |
| | | | | | | | 2-Methylphenol | 10 ug/mL |
| | | | | | | | 2-Nitroaniline | 10 ug/mL |
| | | | | | | | 2-Nitrophenol | 10 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 10 ug/mL |
| | | | | | | | 3-Nitroaniline | 10 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 20 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 10 ug/mL |
| | | | | | | | 4-Chloroaniline | 10 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Nitroaniline | 10 ug/mL |
| | | | | | | | 4-Nitrophenol | 20 ug/mL |
| | | | | | | | Acenaphthene | 10 ug/mL |
| | | | | | | | Acenaphthylene | 10 ug/mL |
| | | | | | | | Acetophenone | 10 ug/mL |
| | | | | | | | Aniline | 10 ug/mL |
| | | | | | | | Anthracene | 10 ug/mL |
| | | | | | | | Azobenzene | 10 ug/mL |
| | | | | | | | Benzo[a]anthracene | 10 ug/mL |
| | | | | | | | Benzo[a]pyrene | 10 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 10 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 10 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 10 ug/mL |
| | | | | | | | Benzyl alcohol | 10 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 10 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 10 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 10 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 10 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 10 ug/mL |
| | | | | | | | Carbazole | 10 ug/mL |
| | | | | | | | Chrysene | 10 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 10 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 10 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 10 ug/mL |
| | | | | | | | Dibenzofuran | 10 ug/mL |
| | | | | | | | Diethyl phthalate | 10 ug/mL |
| | | | | | | | Dimethyl phthalate | 10 ug/mL |
| | | | | | | | Fluoranthene | 10 ug/mL |
| | | | | | | | Fluorene | 10 ug/mL |
| | | | | | | | Hexachlorobenzene | 10 ug/mL |
| | | | | | | | Hexachlorobutadiene | 10 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 10 ug/mL |
| | | | | | | | Hexachloroethane | 10 ug/mL |
| | | | | | | | Hexadecane | 10 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Isophorone | 10 ug/mL |
| | | | | | | | n-Decane | 10 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 10 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 10 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 20 ug/mL |
| | | | | | | | n-Octadecane | 10 ug/mL |
| | | | | | | | Naphthalene | 10 ug/mL |
| | | | | | | | Nitrobenzene | 10 ug/mL |
| | | | | | | | Pentachlorophenol | 20 ug/mL |
| | | | | | | | Phenanthrene | 10 ug/mL |
| | | | | | | | Phenol | 10 ug/mL |
| | | | | | | | Pyrene | 10 ug/mL |
| | | | | | | | Pyridine | 10 ug/mL |
| | | | | | | | Benzoic acid | 20 ug/mL |
| | | | | | | | Indene | 20 ug/mL |
| | | | | | | | Atrazine | 20 ug/mL |
| | | | | | | | Benzaldehyde | 20 ug/mL |
| | | | | | | | Caprolactam | 20 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 20 ug/mL |
| | | | | | | | Benzidine | 20 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 10 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 10 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 10 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 10 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L7 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 300 uL | 1,1'-Biphenyl | 15 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 15 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 15 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 15 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 15 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 15 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 15 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 15 ug/mL |
| | | | | | | | 1,4-Dioxane | 15 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1-Methylnaphthalene | 15 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 15 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 15 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 15 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 15 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 15 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 30 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 15 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 15 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 15 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 15 ug/mL |
| | | | | | | | 2-Chlorophenol | 15 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 15 ug/mL |
| | | | | | | | 2-Methylphenol | 15 ug/mL |
| | | | | | | | 2-Nitroaniline | 15 ug/mL |
| | | | | | | | 2-Nitrophenol | 15 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 15 ug/mL |
| | | | | | | | 3-Nitroaniline | 15 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 30 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 15 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 15 ug/mL |
| | | | | | | | 4-Chloroaniline | 15 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 15 ug/mL |
| | | | | | | | 4-Nitroaniline | 15 ug/mL |
| | | | | | | | 4-Nitrophenol | 30 ug/mL |
| | | | | | | | Acenaphthene | 15 ug/mL |
| | | | | | | | Acenaphthylene | 15 ug/mL |
| | | | | | | | Acetophenone | 15 ug/mL |
| | | | | | | | Aniline | 15 ug/mL |
| | | | | | | | Anthracene | 15 ug/mL |
| | | | | | | | Azobenzene | 15 ug/mL |
| | | | | | | | Benzo[a]anthracene | 15 ug/mL |
| | | | | | | | Benzo[a]pyrene | 15 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 15 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 15 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 15 ug/mL |
| | | | | | | | Benzyl alcohol | 15 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 15 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 15 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 15 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 15 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 15 ug/mL |
| | | | | | | | Carbazole | 15 ug/mL |
| | | | | | | | Chrysene | 15 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 15 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 15 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 15 ug/mL |
| | | | | | | | Dibenzofuran | 15 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Diethyl phthalate | 15 ug/mL |
| | | | | | | | Dimethyl phthalate | 15 ug/mL |
| | | | | | | | Fluoranthene | 15 ug/mL |
| | | | | | | | Fluorene | 15 ug/mL |
| | | | | | | | Hexachlorobenzene | 15 ug/mL |
| | | | | | | | Hexachlorobutadiene | 15 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 15 ug/mL |
| | | | | | | | Hexachloroethane | 15 ug/mL |
| | | | | | | | Hexadecane | 15 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 15 ug/mL |
| | | | | | | | Isophorone | 15 ug/mL |
| | | | | | | | n-Decane | 15 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 15 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 15 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 30 ug/mL |
| | | | | | | | n-Octadecane | 15 ug/mL |
| | | | | | | | Naphthalene | 15 ug/mL |
| | | | | | | | Nitrobenzene | 15 ug/mL |
| | | | | | | | Pentachlorophenol | 30 ug/mL |
| | | | | | | | Phenanthrene | 15 ug/mL |
| | | | | | | | Phenol | 15 ug/mL |
| | | | | | | | Pyrene | 15 ug/mL |
| | | | | | | | Pyridine | 15 ug/mL |
| | | | | | | | Benzoic acid | 30 ug/mL |
| | | | | | | | Indene | 30 ug/mL |
| | | | | | | | Atrazine | 30 ug/mL |
| | | | | | | | Benzaldehyde | 30 ug/mL |
| | | | | | | | Caprolactam | 30 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 30 ug/mL |
| | | | | | | | Benzidine | 30 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 15 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 15 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 15 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 15 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 15 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 15 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L8 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 400 uL | 1,1'-Biphenyl | 20 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 20 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 20 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dioxane | 20 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 20 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 40 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 20 ug/mL |
| | | | | | | | 2-Chlorophenol | 20 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2-Methylphenol | 20 ug/mL |
| | | | | | | | 2-Nitroaniline | 20 ug/mL |
| | | | | | | | 2-Nitrophenol | 20 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 20 ug/mL |
| | | | | | | | 3-Nitroaniline | 20 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 40 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 20 ug/mL |
| | | | | | | | 4-Chloroaniline | 20 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Nitroaniline | 20 ug/mL |
| | | | | | | | 4-Nitrophenol | 40 ug/mL |
| | | | | | | | Acenaphthene | 20 ug/mL |
| | | | | | | | Acenaphthylene | 20 ug/mL |
| | | | | | | | Acetophenone | 20 ug/mL |
| | | | | | | | Aniline | 20 ug/mL |
| | | | | | | | Anthracene | 20 ug/mL |
| | | | | | | | Azobenzene | 20 ug/mL |
| | | | | | | | Benzo[a]anthracene | 20 ug/mL |
| | | | | | | | Benzo[a]pyrene | 20 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 20 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 20 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 20 ug/mL |
| | | | | | | | Benzyl alcohol | 20 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bis(2-chloroethoxy)methane | 20 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 20 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 20 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 20 ug/mL |
| | | | | | | | Carbazole | 20 ug/mL |
| | | | | | | | Chrysene | 20 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 20 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 20 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 20 ug/mL |
| | | | | | | | Dibenzofuran | 20 ug/mL |
| | | | | | | | Diethyl phthalate | 20 ug/mL |
| | | | | | | | Dimethyl phthalate | 20 ug/mL |
| | | | | | | | Fluoranthene | 20 ug/mL |
| | | | | | | | Fluorene | 20 ug/mL |
| | | | | | | | Hexachlorobenzene | 20 ug/mL |
| | | | | | | | Hexachlorobutadiene | 20 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 20 ug/mL |
| | | | | | | | Hexachloroethane | 20 ug/mL |
| | | | | | | | Hexadecane | 20 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 20 ug/mL |
| | | | | | | | Isophorone | 20 ug/mL |
| | | | | | | | n-Decane | 20 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 40 ug/mL |
| | | | | | | | n-Octadecane | 20 ug/mL |
| | | | | | | | Naphthalene | 20 ug/mL |
| | | | | | | | Nitrobenzene | 20 ug/mL |
| | | | | | | | Pentachlorophenol | 40 ug/mL |
| | | | | | | | Phenanthrene | 20 ug/mL |
| | | | | | | | Phenol | 20 ug/mL |
| | | | | | | | Pyrene | 20 ug/mL |
| | | | | | | | Pyridine | 20 ug/mL |
| | | | | | | | Benzoic acid | 40 ug/mL |
| | | | | | | | Indene | 40 ug/mL |
| | | | | | | | Atrazine | 40 ug/mL |
| | | | | | | | Benzaldehyde | 40 ug/mL |
| | | | | | | | Caprolactam | 40 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 40 ug/mL |
| | | | | | | | Benzidine | 40 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 20 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 20 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 20 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 20 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L9 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 500 uL | 1,1'-Biphenyl | 25 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 25 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 25 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 25 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 25 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 25 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 25 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 25 ug/mL |
| | | | | | | | 1,4-Dioxane | 25 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 25 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 25 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 25 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 25 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 25 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 25 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 50 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 25 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 25 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 25 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 25 ug/mL |
| | | | | | | | 2-Chlorophenol | 25 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 25 ug/mL |
| | | | | | | | 2-Methylphenol | 25 ug/mL |
| | | | | | | | 2-Nitroaniline | 25 ug/mL |
| | | | | | | | 2-Nitrophenol | 25 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 25 ug/mL |
| | | | | | | | 3-Nitroaniline | 25 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 50 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 25 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 25 ug/mL |
| | | | | | | | 4-Chloroaniline | 25 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 25 ug/mL |
| | | | | | | | 4-Nitroaniline | 25 ug/mL |
| | | | | | | | 4-Nitrophenol | 50 ug/mL |
| | | | | | | | Acenaphthene | 25 ug/mL |
| | | | | | | | Acenaphthylene | 25 ug/mL |
| | | | | | | | Acetophenone | 25 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Aniline | 25 ug/mL |
| | | | | | | | Anthracene | 25 ug/mL |
| | | | | | | | Azobenzene | 25 ug/mL |
| | | | | | | | Benzo[a]anthracene | 25 ug/mL |
| | | | | | | | Benzo[a]pyrene | 25 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 25 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 25 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 25 ug/mL |
| | | | | | | | Benzyl alcohol | 25 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 25 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 25 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 25 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 25 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 25 ug/mL |
| | | | | | | | Carbazole | 25 ug/mL |
| | | | | | | | Chrysene | 25 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 25 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 25 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 25 ug/mL |
| | | | | | | | Dibenzofuran | 25 ug/mL |
| | | | | | | | Diethyl phthalate | 25 ug/mL |
| | | | | | | | Dimethyl phthalate | 25 ug/mL |
| | | | | | | | Fluoranthene | 25 ug/mL |
| | | | | | | | Fluorene | 25 ug/mL |
| | | | | | | | Hexachlorobenzene | 25 ug/mL |
| | | | | | | | Hexachlorobutadiene | 25 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 25 ug/mL |
| | | | | | | | Hexachloroethane | 25 ug/mL |
| | | | | | | | Hexadecane | 25 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 25 ug/mL |
| | | | | | | | Isophorone | 25 ug/mL |
| | | | | | | | n-Decane | 25 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 25 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 25 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 50 ug/mL |
| | | | | | | | n-Octadecane | 25 ug/mL |
| | | | | | | | Naphthalene | 25 ug/mL |
| | | | | | | | Nitrobenzene | 25 ug/mL |
| | | | | | | | Pentachlorophenol | 50 ug/mL |
| | | | | | | | Phenanthrene | 25 ug/mL |
| | | | | | | | Phenol | 25 ug/mL |
| | | | | | | | Pyrene | 25 ug/mL |
| | | | | | | | Pyridine | 25 ug/mL |
| | | | | | | | Benzoic acid | 50 ug/mL |
| | | | | | | | Indene | 50 ug/mL |
| | | | | | | | Atrazine | 50 ug/mL |
| | | | | | | | Benzaldehyde | 50 ug/mL |
| | | | | | | | Caprolactam | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 3,3'-Dichlorobenzidine | 50 ug/mL |
| | | | | | | | Benzidine | 50 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 25 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 25 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 25 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 25 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 25 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 25 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 SS W_00005** | 04/20/16 | 04/21/15 | MECL2, Lot 0000093871 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| **SMLIST1 SS W_00005** | 04/20/16 | 04/21/15 | MECL2, Lot 0000093871 | 2 mL | SMLIST1 SS ST_00005 | 200 uL | 1,1'-Biphenyl | 10 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 10 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 10 ug/mL |
| | | | | | | | 2-Chlorophenol | 10 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 10 ug/mL |
| | | | | | | | 2-Methylphenol | 10 ug/mL |
| | | | | | | | 2-Nitroaniline | 10 ug/mL |
| | | | | | | | 2-Nitrophenol | 10 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 3-Nitroaniline | 10 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 20 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 10 ug/mL |
| | | | | | | | 4-Chloroaniline | 10 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Nitroaniline | 10 ug/mL |
| | | | | | | | 4-Nitrophenol | 20 ug/mL |
| | | | | | | | Acenaphthene | 10 ug/mL |
| | | | | | | | Acenaphthylene | 10 ug/mL |
| | | | | | | | Acetophenone | 10 ug/mL |
| | | | | | | | Anthracene | 10 ug/mL |
| | | | | | | | Benzo[a]anthracene | 10 ug/mL |
| | | | | | | | Benzo[a]pyrene | 10 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 10 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 10 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 10 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 10 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 10 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 10 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 10 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 10 ug/mL |
| | | | | | | | Carbazole | 10 ug/mL |
| | | | | | | | Chrysene | 10 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 10 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 10 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 10 ug/mL |
| | | | | | | | Dibenzofuran | 10 ug/mL |
| | | | | | | | Diethyl phthalate | 10 ug/mL |
| | | | | | | | Dimethyl phthalate | 10 ug/mL |
| | | | | | | | Fluoranthene | 10 ug/mL |
| | | | | | | | Fluorene | 10 ug/mL |
| | | | | | | | Hexachlorobenzene | 10 ug/mL |
| | | | | | | | Hexachlorobutadiene | 10 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 10 ug/mL |
| | | | | | | | Hexachloroethane | 10 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 10 ug/mL |
| | | | | | | | Isophorone | 10 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 10 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 20 ug/mL |
| | | | | | | | Naphthalene | 10 ug/mL |
| | | | | | | | Nitrobenzene | 10 ug/mL |
| | | | | | | | Pentachlorophenol | 20 ug/mL |
| | | | | | | | Phenanthrene | 10 ug/mL |
| | | | | | | | Phenol | 10 ug/mL |
| | | | | | | | Pyrene | 10 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 20 ug/mL |
| | | | | | | | Atrazine | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzaldehyde | 20 ug/mL |
| | | | | | | | Caprolactam | 20 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 10 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 10 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 10 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 10 ug/mL |
| .SMLIST1 SS ST_00005 | 06/30/16 | 04/21/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 SS S1_00004 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | SMLIST1 SS S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | SMLIST1SS S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1SS SUR_00002 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 SS S1_00004 | 06/30/16 | | Restek, Lot A0107697 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| ..SMLIST1 SS S9_00001 | 07/31/16 | | Restek, Lot A0108170 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| ..SMLIST1SS S11_00001 | 06/30/16 | | Restek, Lot A0108042 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..SMLIST1SS SUR_00002 | 09/30/18 | | Restek, Lot A092712 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **VGGASCALNEW_00014** | 09/22/15 | 06/23/15 | MEOH, Lot 62345 | 25 mL | VGGASSTOCKNEW_00003 | 1 mL | C6-C12 | 100 ug/mL |
| | | | | | | | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGGASSTOCKNEW_00003 | 01/31/20 | | RESTEK, Lot A092645 | | (Purchased Reagent) | | C6-C12 | 2500 ug/mL |
| | | | | | | | Gasoline Range Organics [C6 - C10] | 2500 ug/mL |
| **VGGASCALNEW_00015** | 12/23/15 | 09/23/15 | MEOH, Lot 62345 | 25 mL | VGGASSTOCKNEW_00003 | 1 mL | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGGASSTOCKNEW_00003 | 01/31/20 | | RESTEK, Lot A092645 | | (Purchased Reagent) | | Gasoline Range Organics [C6 - C10] | 2500 ug/mL |
| **VGGASCHK_00068** | 10/27/15 | 07/27/15 | MEOH, Lot 49909 | 50 mL | VGASCHK_00002 | 1 mL | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGASCHK_00002 | 10/30/19 | | AccuStandard, Lot B5040009-01 | | (Purchased Reagent) | | Gasoline Range Organics [C6 - C10] | 5 mg/mL |
| **VGGASCHK_00070** | 02/04/16 | 11/04/15 | MEOH, Lot 49909 | 50 mL | VGASCHK_00002 | 1 mL | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGASCHK_00002 | 10/30/19 | | AccuStandard, Lot B5040009-01 | | (Purchased Reagent) | | Gasoline Range Organics [C6 - C10] | 5 mg/mL |
| **VGGASURROGATE_00034** | 01/09/16 | 10/09/15 | MEOH, Lot 49909 | 50 mL | VGSURRSTKGAS_00019 | 1.25 mL | Trifluorotoluene (Surr) | 50 ug/mL |
| .VGSURRSTKGAS_00019 | 08/31/16 | | ULTRA SCIENTIFIC, Lot CE-1903A | | (Purchased Reagent) | | Trifluorotoluene (Surr) | 2000 ug/mL |
| **VM250-IS_00029** | 10/10/15 | 04/10/15 | MEOH, Lot 0000085233 | 100 mL | vmcus8126_00008 | 1 mL | 1,4-Dichlorobenzene-d4 | 250 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 250 ug/mL |
| | | | | | | | Fluorobenzene | 250 ug/mL |
| .vmcus8126_00008 | 09/30/16 | | Ultra Scientific, Lot CL-3192 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 25000 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 25000 ug/mL |
| | | | | | | | Fluorobenzene | 25000 ug/mL |
| **VM250-IS_00030** | 03/17/16 | 09/17/15 | MEOH, Lot +1632000000099494jl | 100 mL | vmcus8126_00008 | 1 mL | 1,4-Dichlorobenzene-d4 | 250 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 250 ug/mL |
| | | | | | | | Fluorobenzene | 250 ug/mL |
| .vmcus8126_00008 | 09/30/16 | | Ultra Scientific, Lot CL-3192 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 25000 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 25000 ug/mL |
| | | | | | | | Fluorobenzene | 25000 ug/mL |
| **VM250SS_00031** | 11/29/15 | 05/29/15 | MEOH, Lot +161620000085233JB | 50 mL | VM567650_00021 | 5 mL | 1,2-Dichloroethane-d4 (Surr) | 250 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 250 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 250 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 250 ug/mL |
| .VM567650_00021 | 04/30/19 | | Restek, Lot A0102817 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **VM250SS_00033** | 02/09/16 | 11/09/15 | MEOH, Lot +171940000118655JM | 50 mL | VM567650_00023 | 5 mL | 1,2-Dichloroethane-d4 (Surr) | 250 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 250 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 250 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 250 ug/mL |
| .VM567650_00023 | 06/30/19 | | Restek, Lot A0104073 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **VM50IS_00050** | 12/11/15 | 06/11/15 | MEOH, Lot 85233 | 100 mL | vm30241_00002 | 2 mL | 1,4-Dichlorobenzene-d4 | 50 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 50 ug/mL |
| | | | | | | | Fluorobenzene | 50 ug/mL |
| .vm30241_00002 | 11/30/19 | | restek, Lot A0107133 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2500 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 2500 ug/mL |
| | | | | | | | Fluorobenzene | 2500 ug/mL |
| **vm50ss_00213** | 09/02/15 | 08/26/15 | MEOH, Lot na | 2 mL | vm50ss_stk_00065 | 2 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| .vm50ss_stk_00065 | 09/08/15 | 03/08/15 | MEOH, Lot 0000085233 | 200 mL | VM567650_00021 | 4 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| ..VM567650_00021 | 04/30/19 | | Restek, Lot A0102817 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **vm50ss_stk_00066** | 12/11/15 | 06/11/15 | MEOH, Lot 0000085233 | 200 mL | VM567650_00021 | 4 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| .VM567650_00021 | 04/30/19 | | Restek, Lot A0102817 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **VMAROLISTDW_00114** | 09/08/15 | 09/01/15 | MEOH, Lot na | 3 mL | VMACROLSTD_00029 | 3 mL | Acrolein | 250 ug/mL |
| .VMACROLSTD_00029 | 09/30/15 | 09/01/15 | MEOH, Lot 147358 | 20 mL | VM568720_00008 | 250 uL | Acrolein | 250 ug/mL |
| ..VM568720_00008 | 09/30/15 | | restek, Lot A0111006 | | (Purchased Reagent) | | Acrolein | 20000 ug/mL |
| **VMAROLISTDW_00122** | 10/31/15 | 10/25/15 | MEOH, Lot na | 3 mL | VMACROLSTD_00030 | 3 mL | Acrolein | 250 ug/mL |
| .VMACROLSTD_00030 | 10/31/15 | 10/01/15 | MEOH, Lot 0000099494 | 20 mL | VM568720_00009 | 250 uL | Acrolein | 250 ug/mL |
| ..VM568720_00009 | 11/30/15 | | restek, Lot A0112264 | | (Purchased Reagent) | | Acrolein | 20000 ug/mL |
| **VMFASA9W_00076** | 11/04/15 | 10/28/15 | MEOH, Lot NA | 2 mL | VMFASA9_00005 | 2 mL | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |
| .VMFASA9_00005 | 11/30/15 | 06/30/15 | MEOH, Lot +161620000085233jb | 100 mL | VM568723S_00001 | 2.5 mL | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725S_00001 | 2 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| ..VM568723S_00001 | 04/30/16 | restek, Lot A0102749 | | | (Purchased Reagent) | | Diisopropyl ether | 2000 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725S_00001 | 11/30/15 | restek, Lot A0111261 | | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| **VMFASG_00033** | 09/29/15 | 08/29/15 | MEOH, Lot 147358 | 50 mL | VM567645S_00016 | 1.25 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .VM567645S_00016 | 11/30/15 | Restek, Lot A099261 | | | (Purchased Reagent) | | Bromomethane | 2000 ug/mL |
| | | | | | | | Chloroethane | 2000 ug/mL |
| | | | | | | | Chloromethane | 2000 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2000 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2000 ug/mL |
| | | | | | | | Vinyl chloride | 2000 ug/mL |
| **VMFASGW_00120** | 10/29/15 | 10/22/15 | MEOH, Lot NA | 2 mL | VMFASG_00034 | 2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .VMFASG_00034 | 10/29/15 | 09/29/15 | MEOH, Lot 0000099494 | 50 mL | VM567645S_00016 | 1.25 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| ..VM567645S_00016 | 11/30/15 | Restek, Lot A099261 | | | (Purchased Reagent) | | Bromomethane | 2000 ug/mL |
| | | | | | | | Chloroethane | 2000 ug/mL |
| | | | | | | | Chloromethane | 2000 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2000 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2000 ug/mL |
| | | | | | | | Vinyl chloride | 2000 ug/mL |
| **VMFASGW_00122** | 11/11/15 | 11/04/15 | MEOH, Lot NA | 2 mL | VMFASG_00035 | 2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .VMFASG_00035 | 11/30/15 | 10/27/15 | MEOH, Lot 0000118655 | 50 mL | VM567645S_00016 | 1.25 mL | Bromomethane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| ..VM567645S_00016 | 11/30/15 | | Restek, Lot A099261 | | (Purchased Reagent) | | Bromomethane | 2000 ug/mL |
| | | | | | | | Chloroethane | 2000 ug/mL |
| | | | | | | | Chloromethane | 2000 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2000 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2000 ug/mL |
| | | | | | | | Vinyl chloride | 2000 ug/mL |
| **VMFASGW_00123** | 11/19/15 | 11/12/15 | MEOH, Lot NA | 2 mL | VMFASG_00035 | 2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .VMFASG_00035 | 11/30/15 | 10/27/15 | MEOH, Lot 0000118655 | 50 mL | VM567645S_00016 | 1.25 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| ..VM567645S_00016 | 11/30/15 | | Restek, Lot A099261 | | (Purchased Reagent) | | Bromomethane | 2000 ug/mL |
| | | | | | | | Chloroethane | 2000 ug/mL |
| | | | | | | | Chloromethane | 2000 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2000 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2000 ug/mL |
| | | | | | | | Vinyl chloride | 2000 ug/mL |
| **VMFASPW_00114** | 09/08/15 | 09/01/15 | MEOH, Lot n/a | 2 mL | VMRFASP_00022 | 2 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRFASP_00022 | 10/31/15 | 07/28/15 | MEOH, Lot 147358 | 100 mL | VM567641S_00007 | 2.5 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | VM567642S_00005 | 1 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM569723S_00002 | 2 mL | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | VM569724S_00002 | 1 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641S_00007 | 02/29/16 | | Restek, Lot A093733 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| | | | | | | | Xylenes, Total | 4000 ug/mL |
| ..VM567642S_00005 | 02/29/16 | | Restek, Lot A093472 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM569723S_00002 | 03/31/18 | | Restek, Lot A0110013 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2500 ug/mL |
| ..VM569724S_00002 | 10/31/15 | | Restek, Lot A0110474 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMFASPW_00122** | 10/28/15 | 10/21/15 | MEOH, Lot n/a | 2 mL | VMRFASP_00023 | 2 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRFASP_00023 | 10/31/15 | 09/29/15 | MEOH, Lot 99494 | 100 mL | VM567641S_00007 | 2.5 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | VM567642S_00005 | 1 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM569723S_00002 | 2 mL | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | VM569724S_00002 | 1 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641S_00007 | 02/29/16 | | Restek, Lot A093733 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| | | | | | | | Xylenes, Total | 4000 ug/mL |
| ..VM567642S_00005 | 02/29/16 | | Restek, Lot A093472 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM569723S_00002 | 03/31/18 | | Restek, Lot A0110013 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2500 ug/mL |
| ..VM569724S_00002 | 10/31/15 | | Restek, Lot A0110474 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMFASPW_00125** | 11/16/15 | 11/09/15 | MEOH, Lot n/a | 2 mL | VMRFASP_00024 | 2 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRFASP_00024 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 100 mL | VM567641S_00007 | 2.5 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | VM567642S_00005 | 1 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM569723S_00001 | 2 mL | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | VM569724S_00003 | 1 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641S_00007 | 02/29/16 | | Restek, Lot A093733 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| | | | | | | | Xylenes, Total | 4000 ug/mL |
| ..VM567642S_00005 | 02/29/16 | | Restek, Lot A093472 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM569723S_00001 | 01/31/18 | | Restek, Lot A0108188 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2500 ug/mL |
| ..VM569724S_00003 | 01/31/16 | | Restek, Lot A0112364 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRA9W_00119** | 08/17/15 | 08/10/15 | MEOH, Lot NA | 1 mL | VMRA9_00013 | 1 mL | Cyclohexanone | 500 ug/mL |
| | | | | | | | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| .VMRA9_00013 | 11/30/15 | 06/30/15 | MEOH, Lot +161620000085233jb | 50 mL | VM567648_00014 | 1.25 mL | Cyclohexanone | 500 ug/mL |
| | | | | | VM568723_00001 | 1.25 mL | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725_00002 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | vm569726_00001 | 2 mL | 2-Methylnaphthalene | 100 ug/mL |
| ..VM567648_00014 | 02/29/16 | | Restek, Lot A093361 | | (Purchased Reagent) | | Cyclohexanone | 20000 ug/mL |
| ..VM568723_00001 | 12/31/15 | | restek, Lot A099930 | | (Purchased Reagent) | | Acetonitrile | 20000 ug/mL |
| | | | | | | | Diisopropyl ether | 2000 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Propionitrile | 20000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725_00002 | 11/30/15 | | restek, Lot A0111228 | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 2500 ug/mL |
| | | | | | | | 1-Chlorohexane | 2500 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 2500 ug/mL |
| | | | | | | | 2-Nitropropane | 5000 ug/mL |
| | | | | | | | Ethyl acetate | 5000 ug/mL |
| | | | | | | | Ethyl acrylate | 2500 ug/mL |
| | | | | | | | Methacrylonitrile | 25000 ug/mL |
| | | | | | | | Methyl methacrylate | 5000 ug/mL |
| | | | | | | | n-Butanol | 62500 ug/mL |
| | | | | | | | n-Butyl acetate | 2500 ug/mL |
| ..vm569726_00001 | 04/30/17 | | Restek, Lot A0110130 | | (Purchased Reagent) | | 2-Methylnaphthalene | 2500 ug/mL |
| **VMRA9W_00128** | 10/30/15 | 10/23/15 | MEOH, Lot NA | 1 mL | VMRA9_00013 | 1 mL | Cyclohexanone | 500 ug/mL |
| | | | | | | | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| .VMRA9_00013 | 11/30/15 | 06/30/15 | MEOH, Lot +161620000085233jb | 50 mL | VM567648_00014 | 1.25 mL | Cyclohexanone | 500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | VM568723_00001 | 1.25 mL | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725_00002 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | vm569726_00001 | 2 mL | 2-Methylnaphthalene | 100 ug/mL |
| ..VM567648_00014 | 02/29/16 | | Restek, Lot A093361 | | (Purchased Reagent) | | Cyclohexanone | 20000 ug/mL |
| ..VM568723_00001 | 12/31/15 | | restek, Lot A099930 | | (Purchased Reagent) | | Acetonitrile | 20000 ug/mL |
| | | | | | | | Diisopropyl ether | 2000 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Propionitrile | 20000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725_00002 | 11/30/15 | | restek, Lot A0111228 | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 2500 ug/mL |
| | | | | | | | 1-Chlorohexane | 2500 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 2500 ug/mL |
| | | | | | | | 2-Nitropropane | 5000 ug/mL |
| | | | | | | | Ethyl acetate | 5000 ug/mL |
| | | | | | | | Ethyl acrylate | 2500 ug/mL |
| | | | | | | | Methacrylonitrile | 25000 ug/mL |
| | | | | | | | Methyl methacrylate | 5000 ug/mL |
| | | | | | | | n-Butanol | 62500 ug/mL |
| | | | | | | | n-Butyl acetate | 2500 ug/mL |
| ..vm569726_00001 | 04/30/17 | | Restek, Lot A0110130 | | (Purchased Reagent) | | 2-Methylnaphthalene | 2500 ug/mL |
| **VMRA9W_00130** | 11/15/15 | 11/08/15 | MEOH, Lot NA | 1 mL | VMRA9_00013 | 1 mL | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |
| .VMRA9_00013 | 11/30/15 | 06/30/15 | MEOH, Lot +161620000085233jb | 50 mL | VM568723_00001 | 1.25 mL | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725_00002 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| ..VM568723_00001 | 12/31/15 | | restek, Lot A099930 | | (Purchased Reagent) | | Diisopropyl ether | 2000 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725_00002 | 11/30/15 | | restek, Lot A0111228 | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **VMRGAS_00116** | 09/05/15 | 08/29/15 | MEOH, Lot 147358 | 10 mL | vm569722_00002 | 0.2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Butadiene | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Dichlorofluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .vm569722_00002 | 01/31/18 | | Restek, Lot A0108198 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Butadiene | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Dichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMRGAS_00123** | 10/30/15 | 10/23/15 | MEOH, Lot 0000099494 | 10 mL | vm569722_00003 | 0.2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Butadiene | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Dichlorofluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .vm569722_00003 | 04/30/18 | | Restek, Lot A0110070 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Butadiene | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Dichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMRGAS_00125** | 11/12/15 | 11/05/15 | MEOH, Lot 0000118655 | 10 mL | vm569722_00003 | 0.2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .vm569722_00003 | 04/30/18 | | Restek, Lot A0110070 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMRPRIMW_00146** | 09/08/15 | 09/01/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00013 | 1 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRPRIM_00013 | 10/31/15 | 09/01/15 | MEOH, Lot 147358 | 50 mL | VM567641_00013 | 1.25 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | VM567642_00015 | 0.5 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM567643_00020 | 2.5 mL | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | VM569724_00002 | 0.5 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641_00013 | 02/29/16 | | restek, Lot A093581 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,4-Dioxane | 40000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Acrylonitrile | 20000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Cyclohexane | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethyl ether | 2000 ug/mL |
| | | | | | | | Ethyl methacrylate | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Hexane | 2000 ug/mL |
| | | | | | | | Iodomethane | 2000 ug/mL |
| | | | | | | | Isobutyl alcohol | 50000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl acetate | 10000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylcyclohexane | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | n-Heptane | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Tetrahydrofuran | 4000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| ..VM567642_00015 | 02/29/16 | | Restek, Lot A093365 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM567643_00020 | 02/29/16 | | restek, Lot A093368 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2000 ug/mL |
| ..VM569724_00002 | 10/31/15 | | Restek, Lot A0110407 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRPRIMW_00153** | 10/30/15 | 10/23/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00013 | 1 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRPRIM_00013 | 10/31/15 | 09/01/15 | MEOH, Lot 147358 | 50 mL | VM567641_00013 | 1.25 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | VM567642_00015 | 0.5 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM567643_00020 | 2.5 mL | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | VM569724_00002 | 0.5 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641_00013 | 02/29/16 | | restek, Lot A093581 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,4-Dioxane | 40000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Acrylonitrile | 20000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Cyclohexane | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethyl ether | 2000 ug/mL |
| | | | | | | | Ethyl methacrylate | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Hexane | 2000 ug/mL |
| | | | | | | | Iodomethane | 2000 ug/mL |
| | | | | | | | Isobutyl alcohol | 50000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl acetate | 10000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylcyclohexane | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | n-Heptane | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Tetrahydrofuran | 4000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| ..VM567642_00015 | 02/29/16 | | Restek, Lot A093365 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM567643_00020 | 02/29/16 | | restek, Lot A093368 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2000 ug/mL |
| ..VM569724_00002 | 10/31/15 | | Restek, Lot A0110407 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRPRIMW_00156** | 11/16/15 | 11/09/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00014 | 1 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRPRIM_00014 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 50 mL | VM567641_00013 | 1.25 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | VM567642_00015 | 0.5 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM567643_00020 | 2.5 mL | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | VM569724_00003 | 0.5 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641_00013 | 02/29/16 | | restek, Lot A093581 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| | | | | | | | Xylenes, Total | 4000 ug/mL |
| ..VM567642_00015 | 02/29/16 | | Restek, Lot A093365 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM567643_00020 | 02/29/16 | | restek, Lot A093368 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2000 ug/mL |
| ..VM569724_00003 | 01/31/16 | | Restek, Lot A0112975 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |

# Certification Summary

Client: Tetra Tech, Inc. TestAmerica Job ID: 240-57597-1
Project/Site: MRC Soil Remedy Block F UST

| Laboratory | Authority | Program | EPA Region | Certification ID |
|---|---|---|---|---|
| TestAmerica Canton | California | State Program | 9 | 2927 |
| TestAmerica Canton | Connecticut | State Program | 1 | PH-0590 |
| TestAmerica Canton | Illinois | NELAP | 5 | 200004 |
| TestAmerica Canton | Kansas | NELAP | 7 | E-10336 |
| TestAmerica Canton | Kentucky (UST) | State Program | 4 | 58 |
| TestAmerica Canton | Kentucky (WW) | State Program | 4 | 98016 |
| TestAmerica Canton | L-A-B | DoD ELAP | | L2315 |
| TestAmerica Canton | Minnesota | NELAP | 5 | 039-999-348 |
| TestAmerica Canton | Nevada | State Program | 9 | OH-000482008A |
| TestAmerica Canton | New Jersey | NELAP | 2 | OH001 |
| TestAmerica Canton | New York | NELAP | 2 | 10975 |
| TestAmerica Canton | Ohio VAP | State Program | 5 | CL0024 |
| TestAmerica Canton | Oregon | NELAP | 10 | 4062 |
| TestAmerica Canton | Pennsylvania | NELAP | 3 | 68-00340 |
| TestAmerica Canton | Texas | NELAP | 6 | T104704517-15-5 |
| TestAmerica Canton | USDA | Federal | | P330-13-00319 |
| TestAmerica Canton | Virginia | NELAP | 3 | 460175 |
| TestAmerica Canton | Washington | State Program | 10 | C971 |
| TestAmerica Canton | West Virginia DEP | State Program | 3 | 210 |
| TestAmerica Canton | Wisconsin | State Program | 5 | 999518190 |

Accreditation may not be offered or required for all methods and analytes reported in this package. Please contact your project manager for the laboratory's current list of certified methods and analytes.

# Method 8260C

## Volatile Organic Compounds (GC/MS) by Method 8260C

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low

GC Column (1): DB-624  ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | DBFM # | DCA # | TOL # | BFB # |
|---|---|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 101 | 103 | 103 | 91 |
| F-VB-UST-2-14 | 240-57597-2 | 95 | 99 | 96 | 85 |
| F-VB-UST-3-14 | 240-57597-3 | 90 | 90 | 93 | 86 |
| F-VB-UST-4-14 | 240-57597-4 | 94 | 102 | 93 | 83 |
| F-VB-UST-5-14 | 240-57597-5 | 101 | 104 | 100 | 87 |
| F-VB-UST-6-14 | 240-57597-6 | 89 | 93 | 88 | 81 |
| F-VB-UST-7-14 | 240-57597-7 | 98 | 103 | 98 | 87 |
| F-VB-UST-8-14 | 240-57597-8 | 93 | 93 | 97 | 85 |
| F-VB-UST-9-14 | 240-57597-9 | 88 | 94 | 89 | 82 |
| F-VB-UST-10-14 | 240-57597-10 | 88 | 93 | 89 | 83 |
| F-VB-UST-11-14 | 240-57597-11 | 87 | 91 | 90 | 81 |
| F-VB-UST-12-14 | 240-57597-12 | 98 | 103 | 99 | 87 |
| F-VB-UST-13-14 | 240-57597-13 | 98 | 100 | 105 | 90 |
| F-VB-UST-14-14 | 240-57597-14 | 89 | 96 | 87 | 78 |
| F-VB-UST-DUP | 240-57597-15 | 101 | 102 | 102 | 89 |
| | MB 240-206089/1-A | 88 | 91 | 91 | 82 |
| | MB 240-206267/6 | 101 | 102 | 101 | 90 |
| | LCS 240-206051/5 | 95 | 92 | 94 | 87 |
| | LCS 240-206267/5 | 99 | 94 | 95 | 89 |
| F-VB-UST-5-14 MS | 240-57597-5 MS | 97 | 92 | 99 | 95 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 97 | 97 | 97 | 88 |
| F-VB-UST-5-14 MSD | 240-57597-5 MSD | 101 | 98 | 101 | 96 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 99 | 95 | 96 | 86 |

| | QC LIMITS |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 37-132 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 58-123 |
| TOL = Toluene-d8 (Surr) | 67-125 |
| BFB = 4-Bromofluorobenzene (Surr) | 52-136 |

# Column to be used to flag recovery values

FORM II 8260C

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low

GC Column (1): DB-624 ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | DBFM # | DCA # | TOL # | BFB # |
|---|---|---|---|---|---|
| TB-110515 | 240-57597-16 | 97 | 99 | 85 | 90 |
| | MB 240-206461/6 | 99 | 95 | 88 | 97 |
| | LCS 240-206461/4 | 103 | 99 | 96 | 103 |
| | 240-57564-B-1 MS | 99 | 97 | 88 | 103 |
| | 240-57564-B-1 MSD | 101 | 98 | 90 | 102 |

| | QC LIMITS |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 79-120 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 78-125 |
| TOL = Toluene-d8 (Surr) | 80-120 |
| BFB = 4-Bromofluorobenzene (Surr) | 61-120 |

# Column to be used to flag recovery values

FORM II 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146327.D

Lab ID: LCS 240-206051/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 20.4 | 102 | 80-120 | |
| 1,1,1-Trichloroethane | 20.0 | 22.0 | 110 | 77-126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 19.6 | 98 | 77-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 20.9 | 105 | 80-138 | |
| 1,1-Dichloroethane | 20.0 | 18.5 | 93 | 76-120 | |
| 1,1-Dichloroethene | 20.0 | 19.1 | 95 | 75-135 | |
| 1,1-Dichloropropene | 20.0 | 20.3 | 102 | 80-120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 19.5 | 98 | 61-121 | |
| 1,2,3-Trichloropropane | 20.0 | 20.3 | 102 | 80-126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 20.6 | 103 | 64-124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 20.3 | 101 | 80-129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 19.1 | 95 | 61-132 | |
| 1,2-Dichlorobenzene | 20.0 | 19.0 | 95 | 76-120 | |
| 1,2-Dichloroethane | 20.0 | 19.3 | 97 | 72-120 | |
| 1,2-Dichloropropane | 20.0 | 17.7 | 88 | 80-120 | |
| 1,3-Dichlorobenzene | 20.0 | 19.3 | 96 | 78-120 | |
| 1,3-Dichloropropane | 20.0 | 18.6 | 93 | 80-120 | |
| 1,4-Dichlorobenzene | 20.0 | 18.8 | 94 | 75-120 | |
| 2,2-Dichloropropane | 20.0 | 21.3 | 106 | 69-135 | |
| 2-Butanone (MEK) | 40.0 | 31.5 | 79 | 52-131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 14.8 J | 74 | 28-173 | |
| 2-Chlorotoluene | 20.0 | 20.2 | 101 | 78-120 | |
| 2-Hexanone | 40.0 | 31.3 | 78 | 64-136 | |
| 4-Chlorotoluene | 20.0 | 20.0 | 100 | 79-120 | |
| p-Isopropyltoluene | 20.0 | 20.7 | 103 | 77-131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 30.7 | 77 | 67-135 | |
| Acetone | 40.0 | 32.8 | 82 | 41-137 | |
| Benzene | 20.0 | 18.5 | 93 | 79-120 | |
| Bromobenzene | 20.0 | 21.3 | 106 | 80-120 | |
| Bromoform | 20.0 | 16.3 | 82 | 62-133 | |
| Bromomethane | 20.0 | 18.8 | 94 | 42-136 | |
| Carbon disulfide | 20.0 | 15.1 | 75 | 62-146 | |
| Carbon tetrachloride | 20.0 | 21.9 | 110 | 71-129 | |
| Chlorobenzene | 20.0 | 19.9 | 99 | 78-120 | |
| Dibromochloromethane | 20.0 | 18.2 | 91 | 72-127 | |
| Chloroethane | 20.0 | 18.7 | 93 | 58-120 | |
| Chloroform | 20.0 | 19.9 | 100 | 77-120 | |
| Chloromethane | 20.0 | 14.5 | 72 | 50-120 | |
| cis-1,2-Dichloroethene | 20.0 | 19.6 | 98 | 76-120 | |
| cis-1,3-Dichloropropene | 20.0 | 17.3 | 86 | 74-128 | |
| Dibromomethane | 20.0 | 19.1 | 96 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146327.D

Lab ID: LCS 240-206051/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 20.0 | 21.1 | 106 | 79-120 | |
| Dichlorodifluoromethane | 20.0 | 16.4 | 82 | 26-120 | |
| Bromodichloromethane | 20.0 | 18.1 | 90 | 80-122 | |
| tert-Butyl alcohol (TBA) | 200 | 138 J | 69 | 60-125 | |
| Ethylbenzene | 20.0 | 19.8 | 99 | 79-120 | |
| Hexachlorobutadiene | 20.0 | 21.4 | 107 | 54-131 | |
| Isopropylbenzene | 20.0 | 20.6 | 103 | 76-122 | |
| Methyl tert-butyl ether | 20.0 | 18.7 | 93 | 49-165 | |
| Methylene Chloride | 20.0 | 16.7 | 84 | 75-120 | |
| m-Xylene & p-Xylene | 20.0 | 19.8 | 99 | 80-120 | |
| Naphthalene | 20.0 | 18.4 | 92 | 65-123 | |
| n-Butylbenzene | 20.0 | 20.2 | 101 | 68-135 | |
| N-Propylbenzene | 20.0 | 19.8 | 99 | 80-129 | |
| o-Xylene | 20.0 | 20.7 | 103 | 80-120 | |
| sec-Butylbenzene | 20.0 | 19.9 | 100 | 74-129 | |
| Styrene | 20.0 | 18.7 | 93 | 80-120 | |
| Tetrachloroethene | 20.0 | 19.6 | 98 | 79-120 | |
| Toluene | 20.0 | 20.5 | 102 | 75-120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.1 | 100 | 78-120 | |
| trans-1,3-Dichloropropene | 20.0 | 19.3 | 97 | 73-131 | |
| Trichloroethene | 20.0 | 21.1 | 105 | 79-120 | |
| Trichlorofluoromethane | 20.0 | 22.3 | 111 | 57-146 | |
| Vinyl acetate | 20.0 | 16.9 | 85 | 70-130 | |
| Vinyl chloride | 20.0 | 15.8 | 79 | 57-120 | |
| Xylenes, Total | 40.0 | 40.5 | 101 | 80-120 | |
| tert-Butylbenzene | 20.0 | 20.1 | 101 | 76-126 | |
| 1,2-Dibromoethane | 20.0 | 19.7 | 99 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146360.D

Lab ID: LCS 240-206267/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 21.7 | 108 | 80-120 | |
| 1,1,1-Trichloroethane | 20.0 | 22.4 | 112 | 77-126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 20.2 | 101 | 77-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 21.2 | 106 | 80-138 | |
| 1,1-Dichloroethane | 20.0 | 19.0 | 95 | 76-120 | |
| 1,1-Dichloroethene | 20.0 | 18.6 | 93 | 75-135 | |
| 1,1-Dichloropropene | 20.0 | 20.9 | 104 | 80-120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 20.6 | 103 | 61-121 | |
| 1,2,3-Trichloropropane | 20.0 | 21.8 | 109 | 80-126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 21.3 | 106 | 64-124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 21.1 | 106 | 80-129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 20.9 | 104 | 61-132 | |
| 1,2-Dichlorobenzene | 20.0 | 20.6 | 103 | 76-120 | |
| 1,2-Dichloroethane | 20.0 | 20.6 | 103 | 72-120 | |
| 1,2-Dichloropropane | 20.0 | 19.1 | 96 | 80-120 | |
| 1,3-Dichlorobenzene | 20.0 | 20.2 | 101 | 78-120 | |
| 1,3-Dichloropropane | 20.0 | 20.5 | 102 | 80-120 | |
| 1,4-Dichlorobenzene | 20.0 | 20.4 | 102 | 75-120 | |
| 2,2-Dichloropropane | 20.0 | 21.9 | 110 | 69-135 | |
| 2-Butanone (MEK) | 40.0 | 33.7 | 84 | 52-131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 15.3 J | 76 | 28-173 | |
| 2-Chlorotoluene | 20.0 | 21.5 | 108 | 78-120 | |
| 2-Hexanone | 40.0 | 32.9 | 82 | 64-136 | |
| 4-Chlorotoluene | 20.0 | 21.6 | 108 | 79-120 | |
| p-Isopropyltoluene | 20.0 | 22.3 | 112 | 77-131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 32.6 | 82 | 67-135 | |
| Acetone | 40.0 | 33.3 | 83 | 41-137 | |
| Benzene | 20.0 | 19.4 | 97 | 79-120 | |
| Bromobenzene | 20.0 | 22.4 | 112 | 80-120 | |
| Bromoform | 20.0 | 17.7 | 89 | 62-133 | |
| Bromomethane | 20.0 | 19.6 | 98 | 42-136 | |
| Carbon disulfide | 20.0 | 14.8 | 74 | 62-146 | |
| Carbon tetrachloride | 20.0 | 22.8 | 114 | 71-129 | |
| Chlorobenzene | 20.0 | 21.0 | 105 | 78-120 | |
| Dibromochloromethane | 20.0 | 19.4 | 97 | 72-127 | |
| Chloroethane | 20.0 | 19.9 | 99 | 58-120 | |
| Chloroform | 20.0 | 21.0 | 105 | 77-120 | |
| Chloromethane | 20.0 | 15.6 | 78 | 50-120 | |
| cis-1,2-Dichloroethene | 20.0 | 20.1 | 101 | 76-120 | |
| cis-1,3-Dichloropropene | 20.0 | 18.4 | 92 | 74-128 | |
| Dibromomethane | 20.0 | 21.4 | 107 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146360.D

Lab ID: LCS 240-206267/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 20.0 | 22.2 | 111 | 79-120 | |
| Dichlorodifluoromethane | 20.0 | 20.4 | 102 | 26-120 | |
| Bromodichloromethane | 20.0 | 19.2 | 96 | 80-122 | |
| tert-Butyl alcohol (TBA) | 200 | 143 J | 71 | 60-125 | |
| Ethylbenzene | 20.0 | 21.3 | 106 | 79-120 | |
| Hexachlorobutadiene | 20.0 | 23.6 | 118 | 54-131 | |
| Isopropylbenzene | 20.0 | 21.2 | 106 | 76-122 | |
| Methyl tert-butyl ether | 20.0 | 18.5 | 93 | 49-165 | |
| Methylene Chloride | 20.0 | 16.9 | 84 | 75-120 | |
| m-Xylene & p-Xylene | 20.0 | 21.3 | 107 | 80-120 | |
| Naphthalene | 20.0 | 18.4 | 92 | 65-123 | |
| n-Butylbenzene | 20.0 | 21.8 | 109 | 68-135 | |
| N-Propylbenzene | 20.0 | 21.7 | 108 | 80-129 | |
| o-Xylene | 20.0 | 21.7 | 108 | 80-120 | |
| sec-Butylbenzene | 20.0 | 20.9 | 105 | 74-129 | |
| Styrene | 20.0 | 20.0 | 100 | 80-120 | |
| Tetrachloroethene | 20.0 | 21.3 | 107 | 79-120 | |
| Toluene | 20.0 | 21.3 | 106 | 75-120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.3 | 101 | 78-120 | |
| trans-1,3-Dichloropropene | 20.0 | 20.6 | 103 | 73-131 | |
| Trichloroethene | 20.0 | 21.5 | 108 | 79-120 | |
| Trichlorofluoromethane | 20.0 | 22.8 | 114 | 57-146 | |
| Vinyl acetate | 20.0 | 17.7 | 88 | 70-130 | |
| Vinyl chloride | 20.0 | 16.8 | 84 | 57-120 | |
| Xylenes, Total | 40.0 | 43.0 | 108 | 80-120 | |
| tert-Butylbenzene | 20.0 | 21.4 | 107 | 76-126 | |
| 1,2-Dibromoethane | 20.0 | 21.7 | 109 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4393.D

Lab ID: LCS 240-206461/4 Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 10.0 | 11.8 | 118 | 80-120 | |
| 1,1,1-Trichloroethane | 10.0 | 11.1 | 111 | 77-123 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 9.38 | 94 | 71-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 10.0 | 9.57 | 96 | 67-138 | |
| 1,1-Dichloroethane | 10.0 | 9.43 | 94 | 79-125 | |
| 1,1-Dichloroethene | 10.0 | 8.73 | 87 | 76-124 | |
| 1,1-Dichloropropene | 10.0 | 10.4 | 104 | 80-128 | |
| 1,2,3-Trichlorobenzene | 10.0 | 8.64 | 86 | 66-120 | |
| 1,2,3-Trichloropropane | 10.0 | 11.3 | 113 | 72-126 | |
| 1,2,4-Trichlorobenzene | 10.0 | 8.66 | 87 | 61-120 | |
| 1,2,4-Trimethylbenzene | 10.0 | 9.93 | 99 | 76-120 | |
| 1,2-Dibromo-3-Chloropropane | 10.0 | 9.17 | 92 | 50-132 | |
| 1,2-Dichlorobenzene | 10.0 | 9.76 | 98 | 79-120 | |
| 1,2-Dichloroethane | 10.0 | 10.5 | 105 | 80-120 | |
| 1,2-Dichloropropane | 10.0 | 9.07 | 91 | 78-124 | |
| 1,3-Dichlorobenzene | 10.0 | 9.90 | 99 | 79-120 | |
| 1,3-Dichloropropane | 10.0 | 9.29 | 93 | 80-120 | |
| 1,4-Dichlorobenzene | 10.0 | 9.83 | 98 | 79-120 | |
| 2,2-Dichloropropane | 10.0 | 10.4 | 104 | 71-122 | |
| 2-Butanone (MEK) | 20.0 | 16.3 | 82 | 56-138 | |
| 2-Chloroethyl vinyl ether | 10.0 | 8.93 J | 89 | 54-139 | |
| 2-Chlorotoluene | 10.0 | 10.4 | 104 | 80-120 | |
| 2-Hexanone | 20.0 | 17.6 | 88 | 55-141 | |
| 4-Chlorotoluene | 10.0 | 9.89 | 99 | 80-120 | |
| p-Isopropyltoluene | 10.0 | 9.42 | 94 | 77-120 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 18.1 | 90 | 64-135 | |
| Acetone | 20.0 | 13.6 | 68 | 34-148 | |
| Benzene | 10.0 | 9.23 | 92 | 80-120 | |
| Bromobenzene | 10.0 | 10.7 | 107 | 80-120 | |
| Bromoform | 10.0 | 10.7 | 107 | 56-122 | |
| Bromomethane | 10.0 | 11.0 | 110 | 38-132 | |
| Carbon disulfide | 10.0 | 7.86 | 79 | 65-144 | |
| Carbon tetrachloride | 10.0 | 11.6 | 116 | 77-131 | |
| Chlorobenzene | 10.0 | 10.2 | 102 | 80-120 | |
| Dibromochloromethane | 10.0 | 10.6 | 106 | 74-120 | |
| Chloroethane | 10.0 | 9.73 | 97 | 36-126 | |
| Chloroform | 10.0 | 10.6 | 106 | 80-120 | |
| Chloromethane | 10.0 | 8.46 | 85 | 48-133 | |
| cis-1,2-Dichloroethene | 10.0 | 9.74 | 97 | 79-120 | |
| cis-1,3-Dichloropropene | 10.0 | 9.52 | 95 | 74-126 | |
| Dibromomethane | 10.0 | 10.7 | 107 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4393.D

Lab ID: LCS 240-206461/4 Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 10.0 | 10.8 | 108 | 80-120 | |
| Dichlorodifluoromethane | 10.0 | 10.5 | 105 | 23-136 | |
| Bromodichloromethane | 10.0 | 11.0 | 110 | 80-120 | |
| tert-Butyl alcohol (TBA) | 100 | 72.2 | 72 | 38-144 | |
| Ethylbenzene | 10.0 | 10.5 | 105 | 80-120 | |
| Hexachlorobutadiene | 10.0 | 8.37 | 84 | 62-120 | |
| Isopropylbenzene | 10.0 | 10.4 | 104 | 77-120 | |
| Methyl tert-butyl ether | 10.0 | 10.1 | 101 | 69-121 | |
| Methylene Chloride | 10.0 | 9.34 | 93 | 77-129 | |
| m-Xylene & p-Xylene | 10.0 | 10.0 | 100 | 80-120 | |
| Naphthalene | 10.0 | 8.44 | 84 | 53-121 | |
| n-Butylbenzene | 10.0 | 8.71 | 87 | 74-120 | |
| N-Propylbenzene | 10.0 | 10.2 | 102 | 79-123 | |
| o-Xylene | 10.0 | 10.4 | 104 | 80-120 | |
| sec-Butylbenzene | 10.0 | 9.27 | 93 | 77-120 | |
| Styrene | 10.0 | 9.85 | 99 | 76-122 | |
| Tetrachloroethene | 10.0 | 10.9 | 109 | 78-121 | |
| Toluene | 10.0 | 9.77 | 98 | 80-120 | |
| trans-1,2-Dichloroethene | 10.0 | 9.53 | 95 | 80-124 | |
| trans-1,3-Dichloropropene | 10.0 | 9.90 | 99 | 75-131 | |
| Trichloroethene | 10.0 | 11.1 | 111 | 80-121 | |
| Trichlorofluoromethane | 10.0 | 9.70 | 97 | 61-133 | |
| Vinyl acetate | 10.0 | 9.77 | 98 | 27-160 | |
| Vinyl chloride | 10.0 | 9.07 | 91 | 52-121 | |
| Xylenes, Total | 20.0 | 20.4 | 102 | 80-120 | |
| tert-Butylbenzene | 10.0 | 9.67 | 97 | 77-121 | |
| 1,2-Dibromoethane | 10.0 | 10.7 | 107 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146341.D

Lab ID: 240-57597-5 MS Client ID: F-VB-UST-5-14 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 22.8 | 0.58 U | 23.0 | 101 | 34-135 | |
| 1,1,1-Trichloroethane | 22.8 | 0.75 U | 24.1 | 106 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 22.8 | 0.36 U | 26.6 | 117 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 22.8 | 0.49 U | 23.5 | 103 | 50-147 | |
| 1,1-Dichloroethane | 22.8 | 0.39 U | 21.1 | 93 | 54-122 | |
| 1,1-Dichloroethene | 22.8 | 0.93 U | 19.8 | 87 | 49-157 | |
| 1,1-Dichloropropene | 22.8 | 0.43 U | 22.7 | 100 | 50-122 | |
| 1,2,3-Trichlorobenzene | 22.8 | 0.58 U | 14.1 | 62 | 10-120 | |
| 1,2,3-Trichloropropane | 22.8 | 0.90 U | 23.8 | 104 | 32-174 | |
| 1,2,4-Trichlorobenzene | 22.8 | 0.44 U | 14.2 | 62 | 10-120 | |
| 1,2,4-Trimethylbenzene | 22.8 | 0.67 U | 18.7 | 82 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 22.8 | 3.1 U | 22.2 | 97 | 10-153 | |
| 1,2-Dichlorobenzene | 22.8 | 0.39 U | 18.1 | 79 | 17-122 | |
| 1,2-Dichloroethane | 22.8 | 0.53 U | 22.0 | 97 | 49-123 | |
| 1,2-Dichloropropane | 22.8 | 0.16 U | 20.8 | 91 | 61-120 | |
| 1,3-Dichlorobenzene | 22.8 | 0.80 U | 18.3 | 80 | 16-126 | |
| 1,3-Dichloropropane | 22.8 | 0.46 U | 22.5 | 98 | 54-128 | |
| 1,4-Dichlorobenzene | 22.8 | 0.32 U | 17.8 | 78 | 15-121 | |
| 2,2-Dichloropropane | 22.8 | 0.51 U | 24.0 | 105 | 49-132 | |
| 2-Butanone (MEK) | 45.6 | 4.3 J | 36.6 | 71 | 30-143 | |
| 2-Chloroethyl vinyl ether | 22.8 | 0.74 U | 16.4 J | 72 | 10-162 | |
| 2-Chlorotoluene | 22.8 | 0.35 U | 20.8 | 91 | 11-162 | |
| 2-Hexanone | 45.6 | 0.79 U | 52.2 | 114 | 37-147 | |
| 4-Chlorotoluene | 22.8 | 0.30 U | 19.4 | 85 | 15-149 | |
| p-Isopropyltoluene | 22.8 | 0.47 U | 19.0 | 83 | 10-165 | |
| 4-Methyl-2-pentanone (MIBK) | 45.6 | 1.4 U | 150 | 328 | 43-147 | F1 |
| Acetone | 45.6 | 13 J | 47.7 | 77 | 24-140 | |
| Benzene | 22.8 | 1.0 U | 21.1 | 93 | 53-120 | |
| Bromobenzene | 22.8 | 0.38 U | 21.7 | 95 | 24-144 | |
| Bromoform | 22.8 | 0.28 U | 18.7 | 82 | 18-129 | |
| Bromomethane | 22.8 | 0.49 U | 19.8 | 87 | 33-130 | |
| Carbon disulfide | 22.8 | 0.68 U | 16.0 | 70 | 20-151 | |
| Carbon tetrachloride | 22.8 | 0.75 U | 24.0 | 105 | 32-137 | |
| Chlorobenzene | 22.8 | 0.61 U | 20.7 | 91 | 37-120 | |
| Dibromochloromethane | 22.8 | 0.43 U | 20.2 | 89 | 29-135 | |
| Chloroethane | 22.8 | 0.51 U | 21.5 | 94 | 45-120 | |
| Chloroform | 22.8 | 0.42 U | 22.9 | 100 | 53-120 | |
| Chloromethane | 22.8 | 0.92 U | 15.9 | 70 | 34-120 | |
| cis-1,2-Dichloroethene | 22.8 | 0.32 U | 22.1 | 97 | 50-120 | |
| cis-1,3-Dichloropropene | 22.8 | 0.82 U | 19.1 | 84 | 27-133 | |
| Dibromomethane | 22.8 | 0.49 U | 22.0 | 96 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 146341.D

Lab ID: 240-57597-5 MS  Client ID: F-VB-UST-5-14 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Bromochloromethane | 22.8 | 0.38 U | 23.3 | 102 | 53-120 | |
| Dichlorodifluoromethane | 22.8 | 0.32 U | 18.3 | 80 | 17-120 | |
| Bromodichloromethane | 22.8 | 0.34 U | 20.0 | 88 | 35-132 | |
| tert-Butyl alcohol (TBA) | 228 | 6.6 U | 164 J | 72 | 60-125 | |
| Ethylbenzene | 22.8 | 0.37 U | 20.5 | 90 | 30-131 | |
| Hexachlorobutadiene | 22.8 | 0.46 U | 12.5 | 55 | 10-131 | |
| Isopropylbenzene | 22.8 | 0.23 U | 19.5 | 86 | 21-134 | |
| Methyl tert-butyl ether | 22.8 | 0.50 U | 20.9 | 92 | 51-157 | |
| Methylene Chloride | 22.8 | 0.86 U | 18.8 | 82 | 54-120 | |
| m-Xylene & p-Xylene | 22.8 | 0.49 U | 19.3 | 85 | 29-131 | |
| Naphthalene | 22.8 | 0.41 J | 17.8 | 76 | 10-124 | |
| n-Butylbenzene | 22.8 | 0.68 U | 16.6 | 73 | 10-148 | |
| N-Propylbenzene | 22.8 | 0.70 U | 19.7 | 86 | 10-178 | |
| o-Xylene | 22.8 | 0.66 U | 20.9 | 92 | 29-134 | |
| sec-Butylbenzene | 22.8 | 1.1 U | 18.2 | 80 | 10-172 | |
| Styrene | 22.8 | 0.46 U | 17.9 | 79 | 27-127 | |
| Tetrachloroethene | 22.8 | 0.93 U | 22.3 | 98 | 31-135 | |
| Toluene | 22.8 | 0.31 U | 22.0 | 96 | 39-129 | |
| trans-1,2-Dichloroethene | 22.8 | 0.37 U | 22.3 | 98 | 50-123 | |
| trans-1,3-Dichloropropene | 22.8 | 0.34 U | 21.1 | 93 | 28-137 | |
| Trichloroethene | 22.8 | 0.44 U | 23.2 | 101 | 10-177 | |
| Trichlorofluoromethane | 22.8 | 0.34 U | 23.0 | 101 | 36-142 | |
| Vinyl acetate | 22.8 | 0.76 U | 14.4 | 63 | 70-130 | F1 |
| Vinyl chloride | 22.8 | 0.35 U | 17.4 | 76 | 42-120 | |
| Xylenes, Total | 45.6 | 0.63 U | 40.2 | 88 | 30-131 | |
| tert-Butylbenzene | 22.8 | 0.78 U | 21.3 | 93 | 10-163 | |
| 1,2-Dibromoethane | 22.8 | 0.61 U | 24.2 | 106 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146366.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 19.8 | 0.53 U | 20.2 | 102 | 34-135 | |
| 1,1,1-Trichloroethane | 19.8 | 0.68 U | 22.0 | 111 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 19.8 | 0.33 U | 19.2 | 97 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 19.8 | 0.44 U | 21.4 | 108 | 50-147 | |
| 1,1-Dichloroethane | 19.8 | 0.36 U | 18.4 | 93 | 54-122 | |
| 1,1-Dichloroethene | 19.8 | 0.84 U | 18.3 | 93 | 49-157 | |
| 1,1-Dichloropropene | 19.8 | 0.39 U | 20.4 | 103 | 50-122 | |
| 1,2,3-Trichlorobenzene | 19.8 | 0.53 U | 19.1 | 96 | 10-120 | |
| 1,2,3-Trichloropropane | 19.8 | 0.82 U | 22.1 | 112 | 32-174 | |
| 1,2,4-Trichlorobenzene | 19.8 | 0.40 U | 19.5 | 98 | 10-120 | |
| 1,2,4-Trimethylbenzene | 19.8 | 0.61 U | 19.4 | 98 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 19.8 | 2.9 U | 17.8 | 90 | 10-153 | |
| 1,2-Dichlorobenzene | 19.8 | 0.36 U | 18.5 | 94 | 17-122 | |
| 1,2-Dichloroethane | 19.8 | 0.48 U | 19.3 | 97 | 49-123 | |
| 1,2-Dichloropropane | 19.8 | 0.15 U | 17.0 | 86 | 61-120 | |
| 1,3-Dichlorobenzene | 19.8 | 0.73 U | 19.2 | 97 | 16-126 | |
| 1,3-Dichloropropane | 19.8 | 0.42 U | 18.5 | 94 | 54-128 | |
| 1,4-Dichlorobenzene | 19.8 | 0.29 U | 17.9 | 91 | 15-121 | |
| 2,2-Dichloropropane | 19.8 | 0.46 U | 21.6 | 109 | 49-132 | |
| 2-Butanone (MEK) | 39.6 | 3.6 J | 30.1 | 67 | 30-143 | |
| 2-Chloroethyl vinyl ether | 19.8 | 0.67 U | 12.6 J | 63 | 10-162 | |
| 2-Chlorotoluene | 19.8 | 0.32 U | 19.9 | 101 | 11-162 | |
| 2-Hexanone | 39.6 | 0.72 U | 31.1 | 79 | 37-147 | |
| 4-Chlorotoluene | 19.8 | 0.27 U | 19.3 | 98 | 15-149 | |
| p-Isopropyltoluene | 19.8 | 0.43 U | 20.1 | 102 | 10-165 | |
| 4-Methyl-2-pentanone (MIBK) | 39.6 | 1.3 U | 32.3 | 82 | 43-147 | |
| Acetone | 39.6 | 6.0 U | 30.0 | 76 | 24-140 | |
| Benzene | 19.8 | 0.95 U | 18.5 | 94 | 53-120 | |
| Bromobenzene | 19.8 | 0.35 U | 19.9 | 101 | 24-144 | |
| Bromoform | 19.8 | 0.25 U | 16.1 | 81 | 18-129 | |
| Bromomethane | 19.8 | 0.44 U | 19.1 | 97 | 33-130 | |
| Carbon disulfide | 19.8 | 0.62 U | 13.8 | 70 | 20-151 | |
| Carbon tetrachloride | 19.8 | 0.68 U | 21.4 | 108 | 32-137 | |
| Chlorobenzene | 19.8 | 0.56 U | 19.3 | 97 | 37-120 | |
| Dibromochloromethane | 19.8 | 0.39 U | 17.1 | 86 | 29-135 | |
| Chloroethane | 19.8 | 0.46 U | 19.4 | 98 | 45-120 | |
| Chloroform | 19.8 | 0.38 U | 20.5 | 104 | 53-120 | |
| Chloromethane | 19.8 | 0.83 U | 15.5 | 78 | 34-120 | |
| cis-1,2-Dichloroethene | 19.8 | 0.29 U | 19.6 | 99 | 50-120 | |
| cis-1,3-Dichloropropene | 19.8 | 0.75 U | 16.0 | 81 | 27-133 | |
| Dibromomethane | 19.8 | 0.44 U | 21.4 | 108 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 146366.D

Lab ID: 240-57597-15 MS  Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Bromochloromethane | 19.8 | 0.35 U | 20.6 | 104 | 53-120 | |
| Dichlorodifluoromethane | 19.8 | 0.29 U | 19.9 | 101 | 17-120 | |
| Bromodichloromethane | 19.8 | 0.31 U | 17.2 | 87 | 35-132 | |
| tert-Butyl alcohol (TBA) | 198 | 6.0 U | 163 J | 83 | 60-125 | |
| Ethylbenzene | 19.8 | 0.34 U | 19.0 | 96 | 30-131 | |
| Hexachlorobutadiene | 19.8 | 0.42 U | 20.9 | 106 | 10-131 | |
| Isopropylbenzene | 19.8 | 0.21 U | 20.0 | 101 | 21-134 | |
| Methyl tert-butyl ether | 19.8 | 0.45 U | 17.5 | 88 | 51-157 | |
| Methylene Chloride | 19.8 | 0.78 U | 16.2 | 82 | 54-120 | |
| m-Xylene & p-Xylene | 19.8 | 0.44 U | 19.4 | 98 | 29-131 | |
| Naphthalene | 19.8 | 0.37 U | 16.9 | 85 | 10-124 | |
| n-Butylbenzene | 19.8 | 0.62 U | 19.3 | 98 | 10-148 | |
| N-Propylbenzene | 19.8 | 0.63 U | 19.4 | 98 | 10-178 | |
| o-Xylene | 19.8 | 0.60 U | 19.5 | 98 | 29-134 | |
| sec-Butylbenzene | 19.8 | 0.96 U | 19.5 | 98 | 10-172 | |
| Styrene | 19.8 | 0.42 U | 17.8 | 90 | 27-127 | |
| Tetrachloroethene | 19.8 | 0.84 U | 21.7 | 109 | 31-135 | |
| Toluene | 19.8 | 0.28 U | 20.2 | 102 | 39-129 | |
| trans-1,2-Dichloroethene | 19.8 | 0.34 U | 19.7 | 100 | 50-123 | |
| trans-1,3-Dichloropropene | 19.8 | 0.31 U | 17.1 | 86 | 28-137 | |
| Trichloroethene | 19.8 | 0.40 U | 20.3 | 103 | 10-177 | |
| Trichlorofluoromethane | 19.8 | 0.31 U | 21.2 | 107 | 36-142 | |
| Vinyl acetate | 19.8 | 0.70 U | 17.6 | 89 | 70-130 | |
| Vinyl chloride | 19.8 | 0.32 U | 16.9 | 85 | 42-120 | |
| Xylenes, Total | 39.6 | 0.57 U | 38.9 | 98 | 30-131 | |
| tert-Butylbenzene | 19.8 | 0.71 U | 20.0 | 101 | 10-163 | |
| 1,2-Dibromoethane | 19.8 | 0.56 U | 19.8 | 100 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4417.D

Lab ID: 240-57564-B-1 MS Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 2500 | 110 U | 2570 | 103 | 69-122 | |
| 1,1,2,2-Tetrachloroethane | 2500 | 55 U | 2130 | 85 | 61-130 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 2500 | 110 U | 1850 | 74 | 44-140 | |
| 1,1-Dichloroethane | 2500 | 75 U | 2270 | 91 | 73-124 | |
| 1,1-Dichloroethene | 2500 | 110 U | 2020 | 81 | 67-124 | |
| 1,2,4-Trichlorobenzene | 2500 | 80 U | 2090 | 84 | 48-120 | |
| 1,2-Dibromo-3-Chloropropane | 2500 | 210 U | 2310 | 92 | 42-130 | |
| 1,2-Dichlorobenzene | 2500 | 63 U | 2300 | 92 | 67-118 | |
| 1,2-Dichloroethane | 2500 | 58 U | 2580 | 103 | 74-125 | |
| 1,2-Dichloropropane | 2500 | 63 U | 2170 | 87 | 73-122 | |
| 1,3-Dichlorobenzene | 2500 | 48 U | 2250 | 90 | 65-120 | |
| 1,4-Dichlorobenzene | 2500 | 68 U | 2290 | 92 | 66-120 | |
| 2-Butanone (MEK) | 5000 | 7000 | 12400 | 109 | 49-132 | |
| 2-Hexanone | 5000 | 120 U | 4460 | 89 | 49-142 | |
| 4-Methyl-2-pentanone (MIBK) | 5000 | 370 J | 5380 | 100 | 58-136 | |
| Acetone | 5000 | 9800 | 18700 | 179 | 32-126 | F1 |
| Benzene | 2500 | 88 U | 2220 | 89 | 73-121 | |
| Bromoform | 2500 | 140 U | 2500 | 100 | 45-121 | |
| Bromomethane | 2500 | 110 U | 2390 | 96 | 26-136 | |
| Carbon disulfide | 2500 | 95 U | 1870 | 75 | 54-144 | |
| Carbon tetrachloride | 2500 | 110 U | 2620 | 105 | 65-129 | |
| Chlorobenzene | 2500 | 63 U | 2390 | 96 | 72-120 | |
| Dibromochloromethane | 2500 | 110 U | 2460 | 98 | 62-122 | |
| Chloroethane | 2500 | 80 U | 2200 | 88 | 27-131 | |
| Chloroform | 2500 | 63 U | 2560 | 102 | 73-121 | |
| Chloromethane | 2500 | 110 U | 1980 | 79 | 39-134 | |
| cis-1,2-Dichloroethene | 2500 | 65 U | 2370 | 95 | 66-124 | |
| cis-1,3-Dichloropropene | 2500 | 120 U | 2250 | 90 | 60-120 | |
| Dichlorodifluoromethane | 2500 | 80 U | 1870 | 75 | 14-137 | |
| Bromodichloromethane | 2500 | 73 U | 2650 | 106 | 72-120 | |
| Ethylbenzene | 2500 | 63 U | 2440 | 97 | 68-121 | |
| Isopropylbenzene | 2500 | 88 U | 2380 | 95 | 61-122 | |
| Methyl tert-butyl ether | 2500 | 50 U | 2520 | 101 | 61-121 | |
| Methylene Chloride | 2500 | 83 U | 2350 | 94 | 70-124 | |
| m-Xylene & p-Xylene | 2500 | 150 J | 2540 | 96 | 66-123 | |
| o-Xylene | 2500 | 72 J | 2500 | 97 | 68-121 | |
| Styrene | 2500 | 110 U | 2320 | 93 | 64-126 | |
| Tetrachloroethene | 2500 | 78 U | 2450 | 98 | 59-125 | |
| Toluene | 2500 | 170 J | 2380 | 88 | 72-122 | |
| trans-1,2-Dichloroethene | 2500 | 75 U | 2340 | 94 | 72-125 | |
| trans-1,3-Dichloropropene | 2500 | 140 U | 2330 | 93 | 58-132 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4417.D

Lab ID: 240-57564-B-1 MS Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Trichloroethene | 2500 | 55 U | 2580 | 103 | 61-129 | |
| Trichlorofluoromethane | 2500 | 120 U | 2020 | 81 | 49-133 | |
| Vinyl chloride | 2500 | 73 U | 2040 | 81 | 44-122 | |
| Xylenes, Total | 5000 | 220 J | 5040 | 96 | 67-122 | |
| 1,2-Dibromoethane | 2500 | 80 U | 2430 | 97 | 69-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146342.D

Lab ID: 240-57597-5 MSD Client ID: F-VB-UST-5-14 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 23.0 | 24.0 | 104 | 4 | 30 | 34-135 | |
| 1,1,1-Trichloroethane | 23.0 | 25.7 | 112 | 7 | 30 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 23.0 | 28.0 | 121 | 5 | 30 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 23.0 | 25.6 | 111 | 9 | 30 | 50-147 | |
| 1,1-Dichloroethane | 23.0 | 22.5 | 98 | 6 | 30 | 54-122 | |
| 1,1-Dichloroethene | 23.0 | 22.0 | 95 | 10 | 30 | 49-157 | |
| 1,1-Dichloropropene | 23.0 | 25.1 | 109 | 10 | 30 | 50-122 | |
| 1,2,3-Trichlorobenzene | 23.0 | 17.9 | 78 | 24 | 30 | 10-120 | |
| 1,2,3-Trichloropropane | 23.0 | 28.9 | 125 | 19 | 30 | 32-174 | |
| 1,2,4-Trichlorobenzene | 23.0 | 19.9 | 86 | 33 | 30 | 10-120 | F2 |
| 1,2,4-Trimethylbenzene | 23.0 | 25.2 | 109 | 29 | 30 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 23.0 | 26.9 | 117 | 19 | 30 | 10-153 | |
| 1,2-Dichlorobenzene | 23.0 | 22.2 | 96 | 20 | 30 | 17-122 | |
| 1,2-Dichloroethane | 23.0 | 24.1 | 105 | 9 | 30 | 49-123 | |
| 1,2-Dichloropropane | 23.0 | 21.9 | 95 | 5 | 30 | 61-120 | |
| 1,3-Dichlorobenzene | 23.0 | 22.6 | 98 | 21 | 30 | 16-126 | |
| 1,3-Dichloropropane | 23.0 | 23.2 | 101 | 3 | 30 | 54-128 | |
| 1,4-Dichlorobenzene | 23.0 | 22.3 | 97 | 22 | 30 | 15-121 | |
| 2,2-Dichloropropane | 23.0 | 24.5 | 107 | 2 | 30 | 49-132 | |
| 2-Butanone (MEK) | 46.1 | 39.3 | 76 | 7 | 30 | 30-143 | |
| 2-Chloroethyl vinyl ether | 23.0 | 17.2 J | 75 | 5 | 30 | 10-162 | |
| 2-Chlorotoluene | 23.0 | 25.6 | 111 | 21 | 30 | 11-162 | |
| 2-Hexanone | 46.1 | 52.6 | 114 | 1 | 30 | 37-147 | |
| 4-Chlorotoluene | 23.0 | 24.9 | 108 | 25 | 30 | 15-149 | |
| p-Isopropyltoluene | 23.0 | 26.0 | 113 | 31 | 30 | 10-165 | F2 |
| 4-Methyl-2-pentanone (MIBK) | 46.1 | 140 | 303 | 7 | 30 | 43-147 | F1 |
| Acetone | 46.1 | 53.3 | 88 | 11 | 30 | 24-140 | |
| Benzene | 23.0 | 23.0 | 100 | 8 | 30 | 53-120 | |
| Bromobenzene | 23.0 | 26.9 | 117 | 22 | 30 | 24-144 | |
| Bromoform | 23.0 | 19.9 | 86 | 7 | 30 | 18-129 | |
| Bromomethane | 23.0 | 21.9 | 95 | 10 | 30 | 33-130 | |
| Carbon disulfide | 23.0 | 17.6 | 76 | 10 | 30 | 20-151 | |
| Carbon tetrachloride | 23.0 | 26.5 | 115 | 10 | 30 | 32-137 | |
| Chlorobenzene | 23.0 | 23.7 | 103 | 14 | 30 | 37-120 | |
| Dibromochloromethane | 23.0 | 21.2 | 92 | 5 | 30 | 29-135 | |
| Chloroethane | 23.0 | 22.3 | 97 | 4 | 30 | 45-120 | |
| Chloroform | 23.0 | 25.1 | 109 | 9 | 30 | 53-120 | |
| Chloromethane | 23.0 | 17.1 | 74 | 8 | 30 | 34-120 | |
| cis-1,2-Dichloroethene | 23.0 | 24.0 | 104 | 8 | 30 | 50-120 | |
| cis-1,3-Dichloropropene | 23.0 | 21.0 | 91 | 9 | 30 | 27-133 | |
| Dibromomethane | 23.0 | 24.2 | 105 | 9 | 30 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 146342.D

Lab ID: 240-57597-5 MSD  Client ID: F-VB-UST-5-14 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Bromochloromethane | 23.0 | 26.0 | 113 | 11 | 30 | 53-120 | |
| Dichlorodifluoromethane | 23.0 | 18.5 | 80 | 1 | 30 | 17-120 | |
| Bromodichloromethane | 23.0 | 22.9 | 99 | 13 | 30 | 35-132 | |
| tert-Butyl alcohol (TBA) | 230 | 189 J | 82 | 14 | 30 | 60-125 | |
| Ethylbenzene | 23.0 | 24.4 | 106 | 17 | 30 | 30-131 | |
| Hexachlorobutadiene | 23.0 | 22.3 | 97 | 57 | 30 | 10-131 | F2 |
| Isopropylbenzene | 23.0 | 25.1 | 109 | 25 | 30 | 21-134 | |
| Methyl tert-butyl ether | 23.0 | 22.3 | 97 | 6 | 30 | 51-157 | |
| Methylene Chloride | 23.0 | 20.2 | 88 | 7 | 30 | 54-120 | |
| m-Xylene & p-Xylene | 23.0 | 24.0 | 104 | 22 | 30 | 29-131 | |
| Naphthalene | 23.0 | 21.0 | 89 | 16 | 30 | 10-124 | |
| n-Butylbenzene | 23.0 | 25.0 | 109 | 40 | 30 | 10-148 | F2 |
| N-Propylbenzene | 23.0 | 25.5 | 111 | 26 | 30 | 10-178 | |
| o-Xylene | 23.0 | 25.1 | 109 | 18 | 30 | 29-134 | |
| sec-Butylbenzene | 23.0 | 25.1 | 109 | 32 | 30 | 10-172 | F2 |
| Styrene | 23.0 | 21.9 | 95 | 20 | 30 | 27-127 | |
| Tetrachloroethene | 23.0 | 24.7 | 107 | 10 | 30 | 31-135 | |
| Toluene | 23.0 | 24.4 | 106 | 11 | 30 | 39-129 | |
| trans-1,2-Dichloroethene | 23.0 | 24.7 | 107 | 10 | 30 | 50-123 | |
| trans-1,3-Dichloropropene | 23.0 | 22.9 | 99 | 8 | 30 | 28-137 | |
| Trichloroethene | 23.0 | 25.3 | 110 | 9 | 30 | 10-177 | |
| Trichlorofluoromethane | 23.0 | 26.2 | 114 | 13 | 30 | 36-142 | |
| Vinyl acetate | 23.0 | 17.8 | 77 | 21 | 30 | 70-130 | |
| Vinyl chloride | 23.0 | 18.6 | 81 | 7 | 30 | 42-120 | |
| Xylenes, Total | 46.1 | 49.1 | 107 | 20 | 30 | 30-131 | |
| tert-Butylbenzene | 23.0 | 26.7 | 116 | 22 | 30 | 10-163 | |
| 1,2-Dibromoethane | 23.0 | 25.0 | 108 | 3 | 30 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146367.D

Lab ID: 240-57597-15 MSD Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 22.6 | 22.0 | 97 | 9 | 30 | 34-135 | |
| 1,1,1-Trichloroethane | 22.6 | 23.8 | 105 | 8 | 30 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 22.6 | 21.3 | 94 | 11 | 30 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 22.6 | 23.5 | 104 | 10 | 30 | 50-147 | |
| 1,1-Dichloroethane | 22.6 | 20.4 | 90 | 10 | 30 | 54-122 | |
| 1,1-Dichloroethene | 22.6 | 20.3 | 90 | 10 | 30 | 49-157 | |
| 1,1-Dichloropropene | 22.6 | 22.2 | 98 | 8 | 30 | 50-122 | |
| 1,2,3-Trichlorobenzene | 22.6 | 19.9 | 88 | 4 | 30 | 10-120 | |
| 1,2,3-Trichloropropane | 22.6 | 20.5 | 91 | 8 | 30 | 32-174 | |
| 1,2,4-Trichlorobenzene | 22.6 | 20.8 | 92 | 6 | 30 | 10-120 | |
| 1,2,4-Trimethylbenzene | 22.6 | 20.8 | 92 | 7 | 30 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 22.6 | 20.5 | 91 | 14 | 30 | 10-153 | |
| 1,2-Dichlorobenzene | 22.6 | 19.9 | 88 | 7 | 30 | 17-122 | |
| 1,2-Dichloroethane | 22.6 | 21.1 | 93 | 9 | 30 | 49-123 | |
| 1,2-Dichloropropane | 22.6 | 19.2 | 85 | 12 | 30 | 61-120 | |
| 1,3-Dichlorobenzene | 22.6 | 19.0 | 84 | 1 | 30 | 16-126 | |
| 1,3-Dichloropropane | 22.6 | 20.2 | 89 | 9 | 30 | 54-128 | |
| 1,4-Dichlorobenzene | 22.6 | 19.1 | 85 | 7 | 30 | 15-121 | |
| 2,2-Dichloropropane | 22.6 | 24.1 | 107 | 11 | 30 | 49-132 | |
| 2-Butanone (MEK) | 45.2 | 31.6 | 62 | 5 | 30 | 30-143 | |
| 2-Chloroethyl vinyl ether | 22.6 | 14.8 J | 66 | 17 | 30 | 10-162 | |
| 2-Chlorotoluene | 22.6 | 21.8 | 96 | 9 | 30 | 11-162 | |
| 2-Hexanone | 45.2 | 33.3 | 74 | 7 | 30 | 37-147 | |
| 4-Chlorotoluene | 22.6 | 20.1 | 89 | 4 | 30 | 15-149 | |
| p-Isopropyltoluene | 22.6 | 21.4 | 94 | 6 | 30 | 10-165 | |
| 4-Methyl-2-pentanone (MIBK) | 45.2 | 33.2 | 73 | 3 | 30 | 43-147 | |
| Acetone | 45.2 | 29.1 | 64 | 3 | 30 | 24-140 | |
| Benzene | 22.6 | 20.6 | 91 | 11 | 30 | 53-120 | |
| Bromobenzene | 22.6 | 22.2 | 98 | 11 | 30 | 24-144 | |
| Bromoform | 22.6 | 17.5 | 77 | 8 | 30 | 18-129 | |
| Bromomethane | 22.6 | 20.8 | 92 | 8 | 30 | 33-130 | |
| Carbon disulfide | 22.6 | 15.4 | 68 | 11 | 30 | 20-151 | |
| Carbon tetrachloride | 22.6 | 23.7 | 105 | 10 | 30 | 32-137 | |
| Chlorobenzene | 22.6 | 20.8 | 92 | 8 | 30 | 37-120 | |
| Dibromochloromethane | 22.6 | 19.4 | 86 | 13 | 30 | 29-135 | |
| Chloroethane | 22.6 | 20.8 | 92 | 7 | 30 | 45-120 | |
| Chloroform | 22.6 | 22.6 | 100 | 9 | 30 | 53-120 | |
| Chloromethane | 22.6 | 17.2 | 76 | 11 | 30 | 34-120 | |
| cis-1,2-Dichloroethene | 22.6 | 21.5 | 95 | 9 | 30 | 50-120 | |
| cis-1,3-Dichloropropene | 22.6 | 18.1 | 80 | 12 | 30 | 27-133 | |
| Dibromomethane | 22.6 | 22.8 | 101 | 7 | 30 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 146367.D

Lab ID: 240-57597-15 MSD  Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Bromochloromethane | 22.6 | 22.2 | 98 | 8 | 30 | 53-120 | |
| Dichlorodifluoromethane | 22.6 | 21.8 | 96 | 9 | 30 | 17-120 | |
| Bromodichloromethane | 22.6 | 20.4 | 90 | 17 | 30 | 35-132 | |
| tert-Butyl alcohol (TBA) | 226 | 158 J | 70 | 4 | 30 | 60-125 | |
| Ethylbenzene | 22.6 | 20.9 | 93 | 10 | 30 | 30-131 | |
| Hexachlorobutadiene | 22.6 | 22.4 | 99 | 7 | 30 | 10-131 | |
| Isopropylbenzene | 22.6 | 21.8 | 96 | 8 | 30 | 21-134 | |
| Methyl tert-butyl ether | 22.6 | 19.5 | 86 | 11 | 30 | 51-157 | |
| Methylene Chloride | 22.6 | 19.2 | 85 | 17 | 30 | 54-120 | |
| m-Xylene & p-Xylene | 22.6 | 20.2 | 89 | 4 | 30 | 29-131 | |
| Naphthalene | 22.6 | 18.8 | 83 | 11 | 30 | 10-124 | |
| n-Butylbenzene | 22.6 | 20.7 | 92 | 7 | 30 | 10-148 | |
| N-Propylbenzene | 22.6 | 20.7 | 91 | 7 | 30 | 10-178 | |
| o-Xylene | 22.6 | 21.5 | 95 | 10 | 30 | 29-134 | |
| sec-Butylbenzene | 22.6 | 20.6 | 91 | 6 | 30 | 10-172 | |
| Styrene | 22.6 | 19.5 | 86 | 9 | 30 | 27-127 | |
| Tetrachloroethene | 22.6 | 21.7 | 96 | 0 | 30 | 31-135 | |
| Toluene | 22.6 | 21.7 | 96 | 7 | 30 | 39-129 | |
| trans-1,2-Dichloroethene | 22.6 | 21.8 | 97 | 10 | 30 | 50-123 | |
| trans-1,3-Dichloropropene | 22.6 | 19.7 | 87 | 14 | 30 | 28-137 | |
| Trichloroethene | 22.6 | 22.5 | 99 | 10 | 30 | 10-177 | |
| Trichlorofluoromethane | 22.6 | 25.1 | 111 | 17 | 30 | 36-142 | |
| Vinyl acetate | 22.6 | 17.9 | 79 | 2 | 30 | 70-130 | |
| Vinyl chloride | 22.6 | 18.5 | 82 | 9 | 30 | 42-120 | |
| Xylenes, Total | 45.2 | 41.7 | 92 | 7 | 30 | 30-131 | |
| tert-Butylbenzene | 22.6 | 21.5 | 95 | 7 | 30 | 10-163 | |
| 1,2-Dibromoethane | 22.6 | 21.9 | 97 | 10 | 30 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4418.D

Lab ID: 240-57564-B-1 MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 2500 | 2630 | 105 | 2 | 14 | 69-122 | |
| 1,1,2,2-Tetrachloroethane | 2500 | 2220 | 89 | 4 | 18 | 61-130 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 2500 | 2160 | 87 | 15 | 35 | 44-140 | |
| 1,1-Dichloroethane | 2500 | 2310 | 93 | 2 | 14 | 73-124 | |
| 1,1-Dichloroethene | 2500 | 2090 | 84 | 4 | 24 | 67-124 | |
| 1,2,4-Trichlorobenzene | 2500 | 2190 | 88 | 5 | 28 | 48-120 | |
| 1,2-Dibromo-3-Chloropropane | 2500 | 2510 | 100 | 8 | 24 | 42-130 | |
| 1,2-Dichlorobenzene | 2500 | 2370 | 95 | 3 | 15 | 67-118 | |
| 1,2-Dichloroethane | 2500 | 2590 | 104 | 0 | 24 | 74-125 | |
| 1,2-Dichloropropane | 2500 | 2330 | 93 | 7 | 15 | 73-122 | |
| 1,3-Dichlorobenzene | 2500 | 2270 | 91 | 1 | 15 | 65-120 | |
| 1,4-Dichlorobenzene | 2500 | 2330 | 93 | 2 | 16 | 66-120 | |
| 2-Butanone (MEK) | 5000 | 12800 | 116 | 3 | 19 | 49-132 | |
| 2-Hexanone | 5000 | 4330 | 87 | 3 | 27 | 49-142 | |
| 4-Methyl-2-pentanone (MIBK) | 5000 | 5460 | 102 | 2 | 32 | 58-136 | |
| Acetone | 5000 | 18900 | 182 | 1 | 28 | 32-126 | F1 |
| Benzene | 2500 | 2250 | 90 | 1 | 13 | 73-121 | |
| Bromoform | 2500 | 2490 | 99 | 1 | 19 | 45-121 | |
| Bromomethane | 2500 | 2400 | 96 | 0 | 35 | 26-136 | |
| Carbon disulfide | 2500 | 1920 | 77 | 3 | 34 | 54-144 | |
| Carbon tetrachloride | 2500 | 2790 | 111 | 6 | 20 | 65-129 | |
| Chlorobenzene | 2500 | 2400 | 96 | 1 | 15 | 72-120 | |
| Dibromochloromethane | 2500 | 2450 | 98 | 0 | 19 | 62-122 | |
| Chloroethane | 2500 | 2180 | 87 | 1 | 35 | 27-131 | |
| Chloroform | 2500 | 2580 | 103 | 1 | 17 | 73-121 | |
| Chloromethane | 2500 | 2060 | 82 | 4 | 20 | 39-134 | |
| cis-1,2-Dichloroethene | 2500 | 2420 | 97 | 2 | 22 | 66-124 | |
| cis-1,3-Dichloropropene | 2500 | 2330 | 93 | 4 | 21 | 60-120 | |
| Dichlorodifluoromethane | 2500 | 2230 | 89 | 18 | 34 | 14-137 | |
| Bromodichloromethane | 2500 | 2760 | 110 | 4 | 19 | 72-120 | |
| Ethylbenzene | 2500 | 2370 | 95 | 3 | 16 | 68-121 | |
| Isopropylbenzene | 2500 | 2400 | 96 | 1 | 20 | 61-122 | |
| Methyl tert-butyl ether | 2500 | 2570 | 103 | 2 | 12 | 61-121 | |
| Methylene Chloride | 2500 | 2330 | 93 | 1 | 14 | 70-124 | |
| m-Xylene & p-Xylene | 2500 | 2510 | 94 | 2 | 15 | 66-123 | |
| o-Xylene | 2500 | 2490 | 97 | 0 | 14 | 68-121 | |
| Styrene | 2500 | 2340 | 93 | 1 | 15 | 64-126 | |
| Tetrachloroethene | 2500 | 2490 | 100 | 2 | 20 | 59-125 | |
| Toluene | 2500 | 2430 | 91 | 2 | 15 | 72-122 | |
| trans-1,2-Dichloroethene | 2500 | 2400 | 96 | 3 | 25 | 72-125 | |
| trans-1,3-Dichloropropene | 2500 | 2310 | 92 | 1 | 22 | 58-132 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4418.D

Lab ID: 240-57564-B-1 MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Trichloroethene | 2500 | 2730 | 109 | 6 | 14 | 61-129 | |
| Trichlorofluoromethane | 2500 | 2240 | 90 | 10 | 25 | 49-133 | |
| Vinyl chloride | 2500 | 2050 | 82 | 1 | 35 | 44-122 | |
| Xylenes, Total | 5000 | 5000 | 96 | 1 | 14 | 67-122 | |
| 1,2-Dibromoethane | 2500 | 2470 | 99 | 2 | 24 | 69-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 146361.D Lab Sample ID: MB 240-206267/6

Matrix: Solid Heated Purge:(Y/N) Y

Instrument ID: A3UX14 Date Analyzed: 11/11/2015 11:43

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-206267/5 | 146360.D | 11/11/2015 11:22 |
| F-VB-UST-DUP | 240-57597-15 | 146365.D | 11/11/2015 13:09 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 146366.D | 11/11/2015 13:31 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 146367.D | 11/11/2015 13:52 |

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 146331.D Lab Sample ID: MB 240-206089/1-A

Matrix: Solid Heated Purge:(Y/N) Y

Instrument ID: A3UX14 Date Analyzed: 11/10/2015 12:56

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 146336.D | 11/10/2015 14:43 |
| F-VB-UST-2-14 | 240-57597-2 | 146337.D | 11/10/2015 15:05 |
| F-VB-UST-3-14 | 240-57597-3 | 146338.D | 11/10/2015 15:26 |
| F-VB-UST-4-14 | 240-57597-4 | 146339.D | 11/10/2015 15:48 |
| F-VB-UST-5-14 | 240-57597-5 | 146340.D | 11/10/2015 16:09 |
| F-VB-UST-5-14 MS | 240-57597-5 MS | 146341.D | 11/10/2015 16:30 |
| F-VB-UST-5-14 MSD | 240-57597-5 MSD | 146342.D | 11/10/2015 16:52 |
| F-VB-UST-6-14 | 240-57597-6 | 146343.D | 11/10/2015 17:13 |
| F-VB-UST-7-14 | 240-57597-7 | 146344.D | 11/10/2015 17:35 |
| F-VB-UST-8-14 | 240-57597-8 | 146345.D | 11/10/2015 17:56 |
| F-VB-UST-9-14 | 240-57597-9 | 146346.D | 11/10/2015 18:18 |
| F-VB-UST-10-14 | 240-57597-10 | 146347.D | 11/10/2015 18:39 |
| F-VB-UST-11-14 | 240-57597-11 | 146348.D | 11/10/2015 19:01 |
| F-VB-UST-12-14 | 240-57597-12 | 146349.D | 11/10/2015 19:22 |
| F-VB-UST-13-14 | 240-57597-13 | 146350.D | 11/10/2015 19:44 |
| F-VB-UST-14-14 | 240-57597-14 | 146351.D | 11/10/2015 20:05 |

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: UXC4396.D Lab Sample ID: MB 240-206461/6

Matrix: Water Heated Purge:(Y/N) N

Instrument ID: A3UX15 Date Analyzed: 11/12/2015 12:11

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-206461/4 | UXC4393.D | 11/12/2015 11:03 |
| TB-110515 | 240-57597-16 | UXC4397.D | 11/12/2015 12:33 |
| | 240-57564-B-1 MS | UXC4417.D | 11/12/2015 20:05 |
| | 240-57564-B-1 MSD | UXC4418.D | 11/12/2015 20:27 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB14045.D BFB Injection Date: 08/13/2015

Instrument ID: A3UX14 BFB Injection Time: 13:54

Analysis Batch No.: 193150

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 19.1 | |
| 75 | 30.0 - 60.0 % of mass 95 | 45.9 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.7 | |
| 173 | Less than 2.0 % of mass 174 | 0.0 | (0.1)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 70.4 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.1 | (7.3)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 67.7 | (96.1)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.5 | (6.6)2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | IC 240-193150/12 | 144576.D | 08/13/2015 | 17:50 |
| | IC 240-193150/13 | 144577.D | 08/13/2015 | 18:11 |
| | IC 240-193150/14 | 144578.D | 08/13/2015 | 18:33 |
| | IC 240-193150/15 | 144579.D | 08/13/2015 | 18:54 |
| | IC 240-193150/16 | 144580.D | 08/13/2015 | 19:16 |
| | IC 240-193150/17 | 144581.D | 08/13/2015 | 19:37 |
| | IC 240-193150/18 | 144582.D | 08/13/2015 | 19:59 |
| | IC 240-193150/19 | 144583.D | 08/13/2015 | 20:20 |
| | IC 240-193150/20 | 144584.D | 08/13/2015 | 20:42 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB14070.D BFB Injection Date: 09/02/2015

Instrument ID: A3UX14 BFB Injection Time: 13:49

Analysis Batch No.: 196026

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 21.6 | |
| 75 | 30.0 - 60.0 % of mass 95 | 45.0 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 7.0 | |
| 173 | Less than 2.0 % of mass 174 | 0.6 | (0.8)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 70.3 | |
| 175 | 5.0 - 9.0 % of mass 174 | 4.9 | (7.0)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 68.0 | (96.8)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.6 | (6.7)2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | IC 240-196026/3 | 144935.D | 09/02/2015 | 14:32 |
| | IC 240-196026/4 | 144936.D | 09/02/2015 | 14:54 |
| | IC 240-196026/5 | 144937.D | 09/02/2015 | 15:15 |
| | ICIS 240-196026/6 | 144938.D | 09/02/2015 | 15:37 |
| | IC 240-196026/7 | 144939.D | 09/02/2015 | 15:58 |
| | IC 240-196026/8 | 144940.D | 09/02/2015 | 16:20 |
| | IC 240-196026/9 | 144941.D | 09/02/2015 | 16:41 |
| | IC 240-196026/10 | 144942.D | 09/02/2015 | 17:03 |
| | IC 240-196026/11 | 144943.D | 09/02/2015 | 17:24 |
| | ICV 240-196026/13 | 144945.D | 09/02/2015 | 18:07 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB14128.D BFB Injection Date: 11/10/2015

Instrument ID: A3UX14 BFB Injection Time: 10:04

Analysis Batch No.: 206051

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 19.2 | |
| 75 | 30.0 - 60.0 % of mass 95 | 45.9 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.7 | |
| 173 | Less than 2.0 % of mass 174 | 0.0 | (0.0)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 75.8 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.4 | (7.1)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 73.3 | (96.7)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.8 | (6.6)2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-206051/3 | 146325.D | 11/10/2015 | 10:47 |
| | CCV 240-206051/4 | 146326.D | 11/10/2015 | 11:08 |
| | LCS 240-206051/5 | 146327.D | 11/10/2015 | 11:30 |
| | MB 240-206089/1-A | 146331.D | 11/10/2015 | 12:56 |
| F-VB-UST-1-14 | 240-57597-1 | 146336.D | 11/10/2015 | 14:43 |
| F-VB-UST-2-14 | 240-57597-2 | 146337.D | 11/10/2015 | 15:05 |
| F-VB-UST-3-14 | 240-57597-3 | 146338.D | 11/10/2015 | 15:26 |
| F-VB-UST-4-14 | 240-57597-4 | 146339.D | 11/10/2015 | 15:48 |
| F-VB-UST-5-14 | 240-57597-5 | 146340.D | 11/10/2015 | 16:09 |
| F-VB-UST-5-14 MS | 240-57597-5 MS | 146341.D | 11/10/2015 | 16:30 |
| F-VB-UST-5-14 MSD | 240-57597-5 MSD | 146342.D | 11/10/2015 | 16:52 |
| F-VB-UST-6-14 | 240-57597-6 | 146343.D | 11/10/2015 | 17:13 |
| F-VB-UST-7-14 | 240-57597-7 | 146344.D | 11/10/2015 | 17:35 |
| F-VB-UST-8-14 | 240-57597-8 | 146345.D | 11/10/2015 | 17:56 |
| F-VB-UST-9-14 | 240-57597-9 | 146346.D | 11/10/2015 | 18:18 |
| F-VB-UST-10-14 | 240-57597-10 | 146347.D | 11/10/2015 | 18:39 |
| F-VB-UST-11-14 | 240-57597-11 | 146348.D | 11/10/2015 | 19:01 |
| F-VB-UST-12-14 | 240-57597-12 | 146349.D | 11/10/2015 | 19:22 |
| F-VB-UST-13-14 | 240-57597-13 | 146350.D | 11/10/2015 | 19:44 |
| F-VB-UST-14-14 | 240-57597-14 | 146351.D | 11/10/2015 | 20:05 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB14129.D  BFB Injection Date: 11/11/2015

Instrument ID: A3UX14  BFB Injection Time: 09:51

Analysis Batch No.: 206267

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 17.9 | |
| 75 | 30.0 - 60.0 % of mass 95 | 44.2 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.3 | |
| 173 | Less than 2.0 % of mass 174 | 0.0 | (0.0)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 76.6 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.9 | (7.7)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 73.1 | (95.3)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.5 | (6.2)2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-206267/3 | 146358.D | 11/11/2015 | 10:39 |
| | CCV 240-206267/4 | 146359.D | 11/11/2015 | 11:00 |
| | LCS 240-206267/5 | 146360.D | 11/11/2015 | 11:22 |
| | MB 240-206267/6 | 146361.D | 11/11/2015 | 11:43 |
| F-VB-UST-DUP | 240-57597-15 | 146365.D | 11/11/2015 | 13:09 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 146366.D | 11/11/2015 | 13:31 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 146367.D | 11/11/2015 | 13:52 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB460.D BFB Injection Date: 10/28/2015

Instrument ID: A3UX15 BFB Injection Time: 12:37

Analysis Batch No.: 204067

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 17.7 | |
| 75 | 30.0 - 60.0 % of mass 95 | 53.2 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.4 | |
| 173 | Less than 2.0 % of mass 174 | 0.7 | (0.9)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 73.2 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.2 | (7.1)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 70.9 | (96.9)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.6 | (6.5)2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD8260 240-204067/2 | UXC3824A.D | 10/28/2015 | 13:05 |
| | STD8260 240-204067/3 | UXC3825A.D | 10/28/2015 | 13:28 |
| | STD8260 240-204067/4 | UXC3826A.D | 10/28/2015 | 13:51 |
| | STD8260 240-204067/5 | UXC3827A.D | 10/28/2015 | 14:13 |
| | STD8260 240-204067/6 | UXC3828A.D | 10/28/2015 | 14:36 |
| | STD8260 240-204067/7 | UXC3829A.D | 10/28/2015 | 14:59 |
| | ICV 240-204067/14 | UXC3830.D | 10/28/2015 | 15:21 |
| | STD 240-204067/8 | UXC3831.D | 10/28/2015 | 15:44 |
| | STD 240-204067/9 | UXC3832.D | 10/28/2015 | 16:06 |
| | STD 240-204067/10 | UXC3833.D | 10/28/2015 | 16:29 |
| | STD 240-204067/11 | UXC3834.D | 10/28/2015 | 16:52 |
| | STD 240-204067/12 | UXC3835.D | 10/28/2015 | 17:14 |
| | STD 240-204067/13 | UXC3836.D | 10/28/2015 | 17:37 |
| | ICV 240-204067/15 | UXC3837.D | 10/28/2015 | 18:00 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB472.D BFB Injection Date: 11/12/2015

Instrument ID: A3UX15 BFB Injection Time: 10:14

Analysis Batch No.: 206461

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 16.9 | |
| 75 | 30.0 - 60.0 % of mass 95 | 53.2 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.7 | |
| 173 | Less than 2.0 % of mass 174 | 0.6 | (0.8)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 80.0 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.8 | (7.3)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 78.1 | (97.7)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 5.2 | (6.7)2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-206461/2 | UXC4392.D | 11/12/2015 | 10:41 |
| | LCS 240-206461/4 | UXC4393.D | 11/12/2015 | 11:03 |
| | CCV 240-206461/3 | UXC4394.D | 11/12/2015 | 11:26 |
| | MB 240-206461/6 | UXC4396.D | 11/12/2015 | 12:11 |
| TB-110515 | 240-57597-16 | UXC4397.D | 11/12/2015 | 12:33 |
| | 240-57564-B-1 MS | UXC4417.D | 11/12/2015 | 20:05 |
| | 240-57564-B-1 MSD | UXC4418.D | 11/12/2015 | 20:27 |

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: ICIS 240-196026/6  Date Analyzed: 09/02/2015 15:37

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 144938.D  Heated Purge: (Y/N) Y

Calibration ID: 30511

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 1489699 | 6.64 | 910038 | 9.37 | 406298 | 11.34 |
| UPPER LIMIT | | 2979398 | 7.14 | 1820076 | 9.87 | 812596 | 11.84 |
| LOWER LIMIT | | 744850 | 6.14 | 455019 | 8.87 | 203149 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCVIS 240-206051/3 | | 1498068 | 6.64 | 927026 | 9.36 | 430150 | 11.34 |
| CCVIS 240-206267/3 | | 1511907 | 6.64 | 908062 | 9.37 | 416832 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVIS 240-206051/3  Date Analyzed: 11/10/2015 10:47

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 146325.D  Heated Purge: (Y/N) Y

Calibration ID: 30511

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1498068 | 6.64 | 927026 | 9.36 | 430150 | 11.34 |
| UPPER LIMIT | | 2996136 | 7.14 | 1854052 | 9.86 | 860300 | 11.84 |
| LOWER LIMIT | | 749034 | 6.14 | 463513 | 8.86 | 215075 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-206051/4 | | 1428310 | 6.64 | 843706 | 9.36 | 400898 | 11.34 |
| LCS 240-206051/5 | | 1546829 | 6.64 | 933267 | 9.37 | 439911 | 11.34 |
| MB 240-206089/1-A | | 1539926 | 6.64 | 908904 | 9.36 | 440491 | 11.34 |
| 240-57597-1 | F-VB-UST-1-14 | 1339894 | 6.64 | 803741 | 9.36 | 401204 | 11.34 |
| 240-57597-2 | F-VB-UST-2-14 | 1366253 | 6.64 | 812014 | 9.36 | 402470 | 11.34 |
| 240-57597-3 | F-VB-UST-3-14 | 1478321 | 6.64 | 869963 | 9.36 | 422508 | 11.34 |
| 240-57597-4 | F-VB-UST-4-14 | 1379125 | 6.64 | 832253 | 9.36 | 407647 | 11.34 |
| 240-57597-5 | F-VB-UST-5-14 | 1381401 | 6.64 | 834119 | 9.36 | 405215 | 11.34 |
| 240-57597-5 MS | F-VB-UST-5-14 MS | 1327604 | 6.64 | 776997 | 9.36 | 334424 | 11.34 |
| 240-57597-5 MSD | F-VB-UST-5-14 MSD | 1417706 | 6.64 | 855433 | 9.36 | 375351 | 11.34 |
| 240-57597-6 | F-VB-UST-6-14 | 1361610 | 6.64 | 837660 | 9.36 | 414338 | 11.34 |
| 240-57597-7 | F-VB-UST-7-14 | 1397574 | 6.64 | 848234 | 9.36 | 420468 | 11.34 |
| 240-57597-8 | F-VB-UST-8-14 | 1221245 | 6.64 | 712201 | 9.36 | 345698 | 11.34 |
| 240-57597-9 | F-VB-UST-9-14 | 1423543 | 6.64 | 857394 | 9.36 | 412539 | 11.34 |
| 240-57597-10 | F-VB-UST-10-14 | 1539979 | 6.64 | 947956 | 9.36 | 457583 | 11.34 |
| 240-57597-11 | F-VB-UST-11-14 | 1382482 | 6.64 | 822070 | 9.36 | 396271 | 11.34 |
| 240-57597-12 | F-VB-UST-12-14 | 1354784 | 6.64 | 808462 | 9.36 | 408687 | 11.34 |
| 240-57597-13 | F-VB-UST-13-14 | 1299068 | 6.64 | 749701 | 9.36 | 370193 | 11.34 |
| 240-57597-14 | F-VB-UST-14-14 | 1212543 | 6.64 | 728498 | 9.36 | 358599 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVIS 240-206267/3  Date Analyzed: 11/11/2015 10:39

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 146358.D  Heated Purge: (Y/N) Y

Calibration ID: 30511

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1511907 | 6.64 | 908062 | 9.37 | 416832 | 11.34 |
| UPPER LIMIT | | 3023814 | 7.14 | 1816124 | 9.87 | 833664 | 11.84 |
| LOWER LIMIT | | 755954 | 6.14 | 454031 | 8.87 | 208416 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-206267/4 | | 1401932 | 6.64 | 838774 | 9.36 | 393967 | 11.34 |
| LCS 240-206267/5 | | 1488652 | 6.64 | 913203 | 9.36 | 430177 | 11.34 |
| MB 240-206267/6 | | 1376673 | 6.64 | 826304 | 9.37 | 402736 | 11.34 |
| 240-57597-15 | F-VB-UST-DUP | 1258982 | 6.64 | 751722 | 9.36 | 367644 | 11.34 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 953829 | 6.64 | 561900 | 9.36 | 271317 | 11.34 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 1326891 | 6.64 | 791618 | 9.36 | 374548 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: STD8260 240-204067/4  Date Analyzed: 10/28/2015 13:51

Instrument ID: A3UX15  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UXC3826A.D  Heated Purge: (Y/N) N

Calibration ID: 31159

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 2262723 | 4.98 | 1708800 | 7.64 | 900225 | 9.85 |
| UPPER LIMIT | | 4525446 | 5.48 | 3417600 | 8.14 | 1800450 | 10.35 |
| LOWER LIMIT | | 1131362 | 4.48 | 854400 | 7.14 | 450113 | 9.35 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-204067/14 | | 2244648 | 4.98 | 1717477 | 7.64 | 915948 | 9.85 |
| ICV 240-204067/15 | | 2191668 | 4.97 | 1675976 | 7.64 | 861710 | 9.85 |
| CCVIS 240-206461/2 | | 1685909 | 4.97 | 1402013 | 7.64 | 853216 | 9.85 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVIS 240-206461/2  Date Analyzed: 11/12/2015 10:41

Instrument ID: A3UX15  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UXC4392.D  Heated Purge: (Y/N) N

Calibration ID: 31161

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1685909 | 4.97 | 1402013 | 7.64 | 853216 | 9.85 |
| UPPER LIMIT | | 3371818 | 5.47 | 2804026 | 8.14 | 1706432 | 10.35 |
| LOWER LIMIT | | 842955 | 4.47 | 701007 | 7.14 | 426608 | 9.35 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| LCS 240-206461/4 | | 1696051 | 4.97 | 1338674 | 7.64 | 737315 | 9.85 |
| CCV 240-206461/3 | | 1607750 | 4.97 | 1329730 | 7.64 | 731454 | 9.85 |
| MB 240-206461/6 | | 1530091 | 4.97 | 1276242 | 7.64 | 702637 | 9.85 |
| 240-57597-16 | TB-110515 | 1547502 | 4.97 | 1309915 | 7.64 | 691488 | 9.85 |
| 240-57564-B-1 MS | | 1604429 | 4.97 | 1330852 | 7.64 | 805667 | 9.85 |
| 240-57564-B-1 MSD | | 1611108 | 4.97 | 1343144 | 7.64 | 800978 | 9.85 |

FB = Fluorobenzene
FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4
Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 146336.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 5.89(g) | Date Analyzed: 11/10/2015 14:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.81 | U | 5.0 | 0.81 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.79 | U | 5.0 | 0.79 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.91 | U | 5.0 | 0.91 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 146336.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 5.89(g) | Date Analyzed: 11/10/2015 14:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.80 | U | 5.0 | 0.80 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.72 | U | 5.0 | 0.72 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.8 | U | 200 | 5.8 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.75 | U | 5.0 | 0.75 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.35 | U | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.92 | U | 5.0 | 0.92 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.81 | U | 5.0 | 0.81 |
| 108-88-3 | Toluene | 0.27 | U | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 146336.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 5.89(g) | Date Analyzed: 11/10/2015 14:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 103 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 91 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 103 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 146337.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 5.36(g) | Date Analyzed: 11/10/2015 15:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.71 | U | 5.5 | 0.71 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.37 | U | 5.5 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 0.88 | U | 5.5 | 0.88 |
| 563-58-6 | 1,1-Dichloropropene | 0.41 | U | 5.5 | 0.41 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.86 | U | 5.5 | 0.86 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.50 | U | 5.5 | 0.50 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.42 | U | 5.5 | 0.42 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.64 | U | 5.5 | 0.64 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.37 | U | 5.5 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 0.50 | U | 5.5 | 0.50 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.76 | U | 5.5 | 0.76 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.48 | U | 5.5 | 0.48 |
| 78-93-3 | 2-Butanone (MEK) | 4.0 | J B | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.70 | U | 55 | 0.70 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.75 | U | 22 | 0.75 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.5 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.45 | U | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 22 | 1.3 |
| 67-64-1 | Acetone | 6.3 | U | 22 | 6.3 |
| 71-43-2 | Benzene | 0.99 | U | 5.5 | 0.99 |
| 108-86-1 | Bromobenzene | 0.36 | U | 5.5 | 0.36 |
| 75-25-2 | Bromoform | 0.26 | U | 5.5 | 0.26 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.65 | U | 5.5 | 0.65 |
| 56-23-5 | Carbon tetrachloride | 0.71 | U | 5.5 | 0.71 |
| 108-90-7 | Chlorobenzene | 0.58 | U | 5.5 | 0.58 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 146337.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 5.36(g) | Date Analyzed: 11/10/2015 15:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.41 | U | 5.5 | 0.41 |
| 75-00-3 | Chloroethane | 0.48 | U | 5.5 | 0.48 |
| 67-66-3 | Chloroform | 0.39 | U | 5.5 | 0.39 |
| 74-87-3 | Chloromethane | 0.87 | U | 5.5 | 0.87 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.78 | U | 5.5 | 0.78 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.36 | U | 5.5 | 0.36 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.3 | U | 220 | 6.3 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.65 | U | 11 | 0.65 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.47 | U | 5.5 | 0.47 |
| 75-09-2 | Methylene Chloride | 0.81 | U | 5.5 | 0.81 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.38 | U | 5.5 | 0.38 |
| 104-51-8 | n-Butylbenzene | 0.65 | U | 5.5 | 0.65 |
| 103-65-1 | N-Propylbenzene | 0.66 | U | 5.5 | 0.66 |
| 95-47-6 | o-Xylene | 0.63 | U | 5.5 | 0.63 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.5 | 1.0 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.47 | U | 5.5 | 0.47 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.88 | U | 5.5 | 0.88 |
| 108-88-3 | Toluene | 0.30 | U | 5.5 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.42 | U | 5.5 | 0.42 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.72 | U | 11 | 0.72 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.59 | U | 11 | 0.59 |
| 98-06-6 | tert-Butylbenzene | 0.74 | U | 5.5 | 0.74 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 146337.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 5.36(g) | Date Analyzed: 11/10/2015 15:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.58 | U | 5.5 | 0.58 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 96 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 95 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 85 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 99 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 146338.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 5.23(g) | Date Analyzed: 11/10/2015 15:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.76 | U | 5.8 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 5.8 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.91 | U | 5.8 | 0.91 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 5.8 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 5.8 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.48 | U | 5.8 | 0.48 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 6.6 | U | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.69 | U | 5.8 | 0.69 |
| 56-23-5 | Carbon tetrachloride | 0.76 | U | 5.8 | 0.76 |
| 108-90-7 | Chlorobenzene | 0.62 | U | 5.8 | 0.62 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 146338.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 5.23(g) | Date Analyzed: 11/10/2015 15:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 5.8 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.83 | U | 5.8 | 0.83 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 5.8 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.69 | U | 12 | 0.69 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.41 | U | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 0.69 | U | 5.8 | 0.69 |
| 103-65-1 | N-Propylbenzene | 0.70 | U | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.48 | U | 5.8 | 0.48 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.77 | U | 12 | 0.77 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.63 | U | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 0.78 | U | 5.8 | 0.78 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 146338.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 5.23(g) | Date Analyzed: 11/10/2015 15:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.62 | U | 5.8 | 0.62 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 93 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 90 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 86 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 90 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 146339.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 5.43(g) | Date Analyzed: 11/10/2015 15:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.71 | U | 5.5 | 0.71 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.37 | U | 5.5 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 0.87 | U | 5.5 | 0.87 |
| 563-58-6 | 1,1-Dichloropropene | 0.40 | U | 5.5 | 0.40 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.85 | U | 5.5 | 0.85 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.50 | U | 5.5 | 0.50 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.41 | U | 5.5 | 0.41 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.63 | U | 5.5 | 0.63 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.37 | U | 5.5 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 0.50 | U | 5.5 | 0.50 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 5.5 | 0.75 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.48 | U | 5.5 | 0.48 |
| 78-93-3 | 2-Butanone (MEK) | 4.0 | J B | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.70 | U | 55 | 0.70 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.74 | U | 22 | 0.74 |
| 106-43-4 | 4-Chlorotoluene | 0.28 | U | 5.5 | 0.28 |
| 99-87-6 | p-Isopropyltoluene | 0.45 | U | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 22 | 1.3 |
| 67-64-1 | Acetone | 6.2 | U | 22 | 6.2 |
| 71-43-2 | Benzene | 0.98 | U | 5.5 | 0.98 |
| 108-86-1 | Bromobenzene | 0.36 | U | 5.5 | 0.36 |
| 75-25-2 | Bromoform | 0.26 | U | 5.5 | 0.26 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.64 | U | 5.5 | 0.64 |
| 56-23-5 | Carbon tetrachloride | 0.71 | U | 5.5 | 0.71 |
| 108-90-7 | Chlorobenzene | 0.58 | U | 5.5 | 0.58 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 146339.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 5.43(g) | Date Analyzed: 11/10/2015 15:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.40 | U | 5.5 | 0.40 |
| 75-00-3 | Chloroethane | 0.48 | U | 5.5 | 0.48 |
| 67-66-3 | Chloroform | 0.39 | U | 5.5 | 0.39 |
| 74-87-3 | Chloromethane | 0.86 | U | 5.5 | 0.86 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.77 | U | 5.5 | 0.77 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.36 | U | 5.5 | 0.36 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.2 | U | 220 | 6.2 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.64 | U | 11 | 0.64 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.47 | U | 5.5 | 0.47 |
| 75-09-2 | Methylene Chloride | 0.81 | U | 5.5 | 0.81 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.38 | U | 5.5 | 0.38 |
| 104-51-8 | n-Butylbenzene | 0.64 | U | 5.5 | 0.64 |
| 103-65-1 | N-Propylbenzene | 0.65 | U | 5.5 | 0.65 |
| 95-47-6 | o-Xylene | 0.62 | U | 5.5 | 0.62 |
| 135-98-8 | sec-Butylbenzene | 0.99 | U | 5.5 | 0.99 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.47 | U | 5.5 | 0.47 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.87 | U | 5.5 | 0.87 |
| 108-88-3 | Toluene | 0.29 | U | 5.5 | 0.29 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.41 | U | 5.5 | 0.41 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.72 | U | 11 | 0.72 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.59 | U | 11 | 0.59 |
| 98-06-6 | tert-Butylbenzene | 0.73 | U | 5.5 | 0.73 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 146339.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 5.43(g) | Date Analyzed: 11/10/2015 15:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.58 | U | 5.5 | 0.58 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 93 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 94 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 83 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 146340.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.09(g) | Date Analyzed: 11/10/2015 16:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U F2 | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.3 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U F2 | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U F1 | 23 | 1.4 |
| 67-64-1 | Acetone | 13 | J | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 146340.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.09(g) | Date Analyzed: 11/10/2015 16:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U F2 | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.41 | J | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 0.68 | U F2 | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.70 | U | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U F2 | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U F1 | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.63 | U | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 0.78 | U | 5.8 | 0.78 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 146340.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.09(g) | Date Analyzed: 11/10/2015 16:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 100 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 104 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 146343.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.18(g) | Date Analyzed: 11/10/2015 17:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 3.8 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 6.6 | U | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 146343.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.18(g) | Date Analyzed: 11/10/2015 17:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.45 | J | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 146343.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.18(g) | Date Analyzed: 11/10/2015 17:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 88 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 89 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 81 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 146344.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 5.31(g) | Date Analyzed: 11/10/2015 17:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.56 | U | 5.6 | 0.56 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.73 | U | 5.6 | 0.73 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.35 | U | 5.6 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.47 | U | 5.6 | 0.47 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.6 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.90 | U | 5.6 | 0.90 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 5.6 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.56 | U | 5.6 | 0.56 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.88 | U | 5.6 | 0.88 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.52 | U | 5.6 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.43 | U | 5.6 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.65 | U | 5.6 | 0.65 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.6 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.52 | U | 5.6 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.6 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.78 | U | 5.6 | 0.78 |
| 142-28-9 | 1,3-Dichloropropane | 0.45 | U | 5.6 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.6 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.6 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 4.1 | J B | 22 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.72 | U | 56 | 0.72 |
| 95-49-8 | 2-Chlorotoluene | 0.34 | U | 5.6 | 0.34 |
| 591-78-6 | 2-Hexanone | 0.76 | U | 22 | 0.76 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.6 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.46 | U | 5.6 | 0.46 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 6.4 | U | 22 | 6.4 |
| 71-43-2 | Benzene | 1.0 | U | 5.6 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.6 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.6 | 0.27 |
| 74-83-9 | Bromomethane | 0.47 | U | 5.6 | 0.47 |
| 75-15-0 | Carbon disulfide | 0.66 | U | 5.6 | 0.66 |
| 56-23-5 | Carbon tetrachloride | 0.73 | U | 5.6 | 0.73 |
| 108-90-7 | Chlorobenzene | 0.60 | U | 5.6 | 0.60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 146344.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 5.31(g) | Date Analyzed: 11/10/2015 17:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.42 | U | 5.6 | 0.42 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.6 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.6 | 0.40 |
| 74-87-3 | Chloromethane | 0.89 | U | 5.6 | 0.89 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.6 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.80 | U | 5.6 | 0.80 |
| 74-95-3 | Dibromomethane | 0.47 | U | 5.6 | 0.47 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.6 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.6 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.6 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.4 | U | 220 | 6.4 |
| 100-41-4 | Ethylbenzene | 0.36 | U | 5.6 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 0.45 | U | 5.6 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.66 | U | 11 | 0.66 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.6 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.6 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.83 | U | 5.6 | 0.83 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.47 | U | 11 | 0.47 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.6 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.66 | U | 5.6 | 0.66 |
| 103-65-1 | N-Propylbenzene | 0.67 | U | 5.6 | 0.67 |
| 95-47-6 | o-Xylene | 0.64 | U | 5.6 | 0.64 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.6 | 1.0 |
| 100-42-5 | Styrene | 0.45 | U | 5.6 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.6 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.46 | U | 5.6 | 0.46 |
| 127-18-4 | Tetrachloroethene | 0.90 | U | 5.6 | 0.90 |
| 108-88-3 | Toluene | 0.30 | U | 5.6 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.36 | U | 5.6 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.6 | 0.33 |
| 79-01-6 | Trichloroethene | 0.43 | U | 5.6 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.6 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.74 | U | 11 | 0.74 |
| 75-01-4 | Vinyl chloride | 0.34 | U | 5.6 | 0.34 |
| 1330-20-7 | Xylenes, Total | 0.61 | U | 11 | 0.61 |
| 98-06-6 | tert-Butylbenzene | 0.75 | U | 5.6 | 0.75 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 146344.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 5.31(g) | Date Analyzed: 11/10/2015 17:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.60 | U | 5.6 | 0.60 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 98 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 103 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 146345.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 5.29(g) | Date Analyzed: 11/10/2015 17:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.48 | U | 5.8 | 0.48 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.92 | U | 5.8 | 0.92 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.3 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.78 | U | 23 | 0.78 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 6.6 | U | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.48 | U | 5.8 | 0.48 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 146345.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 5.29(g) | Date Analyzed: 11/10/2015 17:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.41 | U | 5.8 | 0.41 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.48 | U | 5.8 | 0.48 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.8 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.85 | U | 5.8 | 0.85 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.48 | U | 12 | 0.48 |
| 91-20-3 | Naphthalene | 0.40 | U | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.8 | 1.0 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.92 | U | 5.8 | 0.92 |
| 108-88-3 | Toluene | 0.33 | J | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.8 | 0.33 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.8 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 146345.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 5.29(g) | Date Analyzed: 11/10/2015 17:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 97 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 93 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 85 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 146346.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.35(g) | Date Analyzed: 11/10/2015 18:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.92 | U | 5.8 | 0.92 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.4 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 12 | J | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 146346.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.35(g) | Date Analyzed: 11/10/2015 18:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.85 | U | 5.8 | 0.85 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.40 | U | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.92 | U | 5.8 | 0.92 |
| 108-88-3 | Toluene | 0.36 | J | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 146346.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.35(g) | Date Analyzed: 11/10/2015 18:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 88 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 82 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 146347.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.14(g) | Date Analyzed: 11/10/2015 18:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.56 | U | 5.6 | 0.56 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.73 | U | 5.6 | 0.73 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.35 | U | 5.6 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.47 | U | 5.6 | 0.47 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.6 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.90 | U | 5.6 | 0.90 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 5.6 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.56 | U | 5.6 | 0.56 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.88 | U | 5.6 | 0.88 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.52 | U | 5.6 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.43 | U | 5.6 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.65 | U | 5.6 | 0.65 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.6 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.52 | U | 5.6 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.6 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.78 | U | 5.6 | 0.78 |
| 142-28-9 | 1,3-Dichloropropane | 0.45 | U | 5.6 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.6 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.6 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 3.5 | J B | 22 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.72 | U | 56 | 0.72 |
| 95-49-8 | 2-Chlorotoluene | 0.34 | U | 5.6 | 0.34 |
| 591-78-6 | 2-Hexanone | 0.76 | U | 22 | 0.76 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.6 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.46 | U | 5.6 | 0.46 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 6.4 | U | 22 | 6.4 |
| 71-43-2 | Benzene | 1.0 | U | 5.6 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.6 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.6 | 0.27 |
| 74-83-9 | Bromomethane | 0.47 | U | 5.6 | 0.47 |
| 75-15-0 | Carbon disulfide | 0.66 | U | 5.6 | 0.66 |
| 56-23-5 | Carbon tetrachloride | 0.73 | U | 5.6 | 0.73 |
| 108-90-7 | Chlorobenzene | 0.60 | U | 5.6 | 0.60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 146347.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.14(g) | Date Analyzed: 11/10/2015 18:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.42 | U | 5.6 | 0.42 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.6 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.6 | 0.40 |
| 74-87-3 | Chloromethane | 0.89 | U | 5.6 | 0.89 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.6 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.80 | U | 5.6 | 0.80 |
| 74-95-3 | Dibromomethane | 0.47 | U | 5.6 | 0.47 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.6 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.6 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.6 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.4 | U | 220 | 6.4 |
| 100-41-4 | Ethylbenzene | 0.36 | U | 5.6 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 0.45 | U | 5.6 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.66 | U | 11 | 0.66 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.6 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.6 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.83 | U | 5.6 | 0.83 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.47 | U | 11 | 0.47 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.6 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.66 | U | 5.6 | 0.66 |
| 103-65-1 | N-Propylbenzene | 0.67 | U | 5.6 | 0.67 |
| 95-47-6 | o-Xylene | 0.64 | U | 5.6 | 0.64 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.6 | 1.0 |
| 100-42-5 | Styrene | 0.45 | U | 5.6 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.6 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.46 | U | 5.6 | 0.46 |
| 127-18-4 | Tetrachloroethene | 0.90 | U | 5.6 | 0.90 |
| 108-88-3 | Toluene | 0.30 | U | 5.6 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.36 | U | 5.6 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.6 | 0.33 |
| 79-01-6 | Trichloroethene | 0.43 | U | 5.6 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.6 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.74 | U | 11 | 0.74 |
| 75-01-4 | Vinyl chloride | 0.34 | U | 5.6 | 0.34 |
| 1330-20-7 | Xylenes, Total | 0.61 | U | 11 | 0.61 |
| 98-06-6 | tert-Butylbenzene | 0.75 | U | 5.6 | 0.75 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 146347.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.14(g) | Date Analyzed: 11/10/2015 18:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.60 | U | 5.6 | 0.60 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 88 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 83 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 146348.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/10/2015 19:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.60 | U | 6.0 | 0.60 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.77 | U | 6.0 | 0.77 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.37 | U | 6.0 | 0.37 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.50 | U | 6.0 | 0.50 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 6.0 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.95 | U | 6.0 | 0.95 |
| 563-58-6 | 1,1-Dichloropropene | 0.44 | U | 6.0 | 0.44 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.60 | U | 6.0 | 0.60 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.93 | U | 6.0 | 0.93 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.55 | U | 6.0 | 0.55 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.45 | U | 6.0 | 0.45 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.69 | U | 6.0 | 0.69 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 6.0 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.55 | U | 6.0 | 0.55 |
| 78-87-5 | 1,2-Dichloropropane | 0.17 | U | 6.0 | 0.17 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.82 | U | 6.0 | 0.82 |
| 142-28-9 | 1,3-Dichloropropane | 0.48 | U | 6.0 | 0.48 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 6.0 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.52 | U | 6.0 | 0.52 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 24 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.76 | U | 60 | 0.76 |
| 95-49-8 | 2-Chlorotoluene | 0.36 | U | 6.0 | 0.36 |
| 591-78-6 | 2-Hexanone | 0.81 | U | 24 | 0.81 |
| 106-43-4 | 4-Chlorotoluene | 0.31 | U | 6.0 | 0.31 |
| 99-87-6 | p-Isopropyltoluene | 0.49 | U | 6.0 | 0.49 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.5 | U | 24 | 1.5 |
| 67-64-1 | Acetone | 6.8 | U | 24 | 6.8 |
| 71-43-2 | Benzene | 1.1 | U | 6.0 | 1.1 |
| 108-86-1 | Bromobenzene | 0.39 | U | 6.0 | 0.39 |
| 75-25-2 | Bromoform | 0.29 | U | 6.0 | 0.29 |
| 74-83-9 | Bromomethane | 0.50 | U | 6.0 | 0.50 |
| 75-15-0 | Carbon disulfide | 0.70 | U | 6.0 | 0.70 |
| 56-23-5 | Carbon tetrachloride | 0.77 | U | 6.0 | 0.77 |
| 108-90-7 | Chlorobenzene | 0.63 | U | 6.0 | 0.63 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 146348.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/10/2015 19:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.44 | U | 6.0 | 0.44 |
| 75-00-3 | Chloroethane | 0.52 | U | 6.0 | 0.52 |
| 67-66-3 | Chloroform | 0.43 | U | 6.0 | 0.43 |
| 74-87-3 | Chloromethane | 0.94 | U | 6.0 | 0.94 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 6.0 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.85 | U | 6.0 | 0.85 |
| 74-95-3 | Dibromomethane | 0.50 | U | 6.0 | 0.50 |
| 74-97-5 | Bromochloromethane | 0.39 | U | 6.0 | 0.39 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 6.0 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.35 | U | 6.0 | 0.35 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.8 | U | 240 | 6.8 |
| 100-41-4 | Ethylbenzene | 0.38 | U | 6.0 | 0.38 |
| 87-68-3 | Hexachlorobutadiene | 0.48 | U | 6.0 | 0.48 |
| 108-20-3 | Diisopropyl ether | 0.70 | U | 12 | 0.70 |
| 98-82-8 | Isopropylbenzene | 0.24 | U | 6.0 | 0.24 |
| 1634-04-4 | Methyl tert-butyl ether | 0.51 | U | 6.0 | 0.51 |
| 75-09-2 | Methylene Chloride | 0.88 | U | 6.0 | 0.88 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.50 | U | 12 | 0.50 |
| 91-20-3 | Naphthalene | 0.42 | U | 6.0 | 0.42 |
| 104-51-8 | n-Butylbenzene | 0.70 | U | 6.0 | 0.70 |
| 103-65-1 | N-Propylbenzene | 0.71 | U | 6.0 | 0.71 |
| 95-47-6 | o-Xylene | 0.68 | U | 6.0 | 0.68 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 6.0 | 1.1 |
| 100-42-5 | Styrene | 0.48 | U | 6.0 | 0.48 |
| 994-05-8 | Tert-amyl methyl ether | 0.51 | U | 6.0 | 0.51 |
| 637-92-3 | Ethyl t-butyl ether | 0.49 | U | 6.0 | 0.49 |
| 127-18-4 | Tetrachloroethene | 0.95 | U | 6.0 | 0.95 |
| 108-88-3 | Toluene | 0.32 | U | 6.0 | 0.32 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.38 | U | 6.0 | 0.38 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.35 | U | 6.0 | 0.35 |
| 79-01-6 | Trichloroethene | 0.45 | U | 6.0 | 0.45 |
| 75-69-4 | Trichlorofluoromethane | 0.35 | U | 6.0 | 0.35 |
| 108-05-4 | Vinyl acetate | 0.79 | U | 12 | 0.79 |
| 75-01-4 | Vinyl chloride | 0.36 | U | 6.0 | 0.36 |
| 1330-20-7 | Xylenes, Total | 0.64 | U | 12 | 0.64 |
| 98-06-6 | tert-Butylbenzene | 0.80 | U | 6.0 | 0.80 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 146348.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/10/2015 19:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.63 | U | 6.0 | 0.63 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 90 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 87 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 81 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 91 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 146349.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.37(g) | Date Analyzed: 11/10/2015 19:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.72 | U | 5.5 | 0.72 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.5 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.89 | U | 5.5 | 0.89 |
| 563-58-6 | 1,1-Dichloropropene | 0.41 | U | 5.5 | 0.41 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.86 | U | 5.5 | 0.86 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.51 | U | 5.5 | 0.51 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.42 | U | 5.5 | 0.42 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.64 | U | 5.5 | 0.64 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.5 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.51 | U | 5.5 | 0.51 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.76 | U | 5.5 | 0.76 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.5 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 3.7 | J B | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.71 | U | 55 | 0.71 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.75 | U | 22 | 0.75 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.5 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.45 | U | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 6.3 | U | 22 | 6.3 |
| 71-43-2 | Benzene | 1.0 | U | 5.5 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.5 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.5 | 0.27 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.65 | U | 5.5 | 0.65 |
| 56-23-5 | Carbon tetrachloride | 0.72 | U | 5.5 | 0.72 |
| 108-90-7 | Chlorobenzene | 0.59 | U | 5.5 | 0.59 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 146349.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.37(g) | Date Analyzed: 11/10/2015 19:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.41 | U | 5.5 | 0.41 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.5 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.5 | 0.40 |
| 74-87-3 | Chloromethane | 0.87 | U | 5.5 | 0.87 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.79 | U | 5.5 | 0.79 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.5 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.3 | U | 220 | 6.3 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.65 | U | 11 | 0.65 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.5 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.82 | U | 5.5 | 0.82 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.5 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.65 | U | 5.5 | 0.65 |
| 103-65-1 | N-Propylbenzene | 0.66 | U | 5.5 | 0.66 |
| 95-47-6 | o-Xylene | 0.63 | U | 5.5 | 0.63 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.5 | 1.0 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.5 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.89 | U | 5.5 | 0.89 |
| 108-88-3 | Toluene | 0.30 | J | 5.5 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.42 | U | 5.5 | 0.42 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.73 | U | 11 | 0.73 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.60 | U | 11 | 0.60 |
| 98-06-6 | tert-Butylbenzene | 0.74 | U | 5.5 | 0.74 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 146349.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.37(g) | Date Analyzed: 11/10/2015 19:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.59 | U | 5.5 | 0.59 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 99 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 103 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 146350.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 5.67(g) | Date Analyzed: 11/10/2015 19:44

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.52 | U | 5.2 | 0.52 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.67 | U | 5.2 | 0.67 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.32 | U | 5.2 | 0.32 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.43 | U | 5.2 | 0.43 |
| 75-34-3 | 1,1-Dichloroethane | 0.35 | U | 5.2 | 0.35 |
| 75-35-4 | 1,1-Dichloroethene | 0.83 | U | 5.2 | 0.83 |
| 563-58-6 | 1,1-Dichloropropene | 0.38 | U | 5.2 | 0.38 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.52 | U | 5.2 | 0.52 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.81 | U | 5.2 | 0.81 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.2 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.39 | U | 5.2 | 0.39 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.60 | U | 5.2 | 0.60 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.8 | U | 10 | 2.8 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.35 | U | 5.2 | 0.35 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.2 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.2 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.71 | U | 5.2 | 0.71 |
| 142-28-9 | 1,3-Dichloropropane | 0.41 | U | 5.2 | 0.41 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.29 | U | 5.2 | 0.29 |
| 594-20-7 | 2,2-Dichloropropane | 0.45 | U | 5.2 | 0.45 |
| 78-93-3 | 2-Butanone (MEK) | 3.1 | J B | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.66 | U | 52 | 0.66 |
| 95-49-8 | 2-Chlorotoluene | 0.31 | U | 5.2 | 0.31 |
| 591-78-6 | 2-Hexanone | 0.70 | U | 21 | 0.70 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.2 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.42 | U | 5.2 | 0.42 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 5.9 | U | 21 | 5.9 |
| 71-43-2 | Benzene | 0.93 | U | 5.2 | 0.93 |
| 108-86-1 | Bromobenzene | 0.34 | U | 5.2 | 0.34 |
| 75-25-2 | Bromoform | 0.25 | U | 5.2 | 0.25 |
| 74-83-9 | Bromomethane | 0.43 | U | 5.2 | 0.43 |
| 75-15-0 | Carbon disulfide | 0.61 | U | 5.2 | 0.61 |
| 56-23-5 | Carbon tetrachloride | 0.67 | U | 5.2 | 0.67 |
| 108-90-7 | Chlorobenzene | 0.55 | U | 5.2 | 0.55 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 146350.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 5.67(g) | Date Analyzed: 11/10/2015 19:44

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.38 | U | 5.2 | 0.38 |
| 75-00-3 | Chloroethane | 0.45 | U | 5.2 | 0.45 |
| 67-66-3 | Chloroform | 0.37 | U | 5.2 | 0.37 |
| 74-87-3 | Chloromethane | 0.82 | U | 5.2 | 0.82 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.29 | U | 5.2 | 0.29 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.73 | U | 5.2 | 0.73 |
| 74-95-3 | Dibromomethane | 0.43 | U | 5.2 | 0.43 |
| 74-97-5 | Bromochloromethane | 0.34 | U | 5.2 | 0.34 |
| 75-71-8 | Dichlorodifluoromethane | 0.29 | U | 5.2 | 0.29 |
| 75-27-4 | Bromodichloromethane | 0.30 | U | 5.2 | 0.30 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.9 | U | 210 | 5.9 |
| 100-41-4 | Ethylbenzene | 0.33 | U | 5.2 | 0.33 |
| 87-68-3 | Hexachlorobutadiene | 0.41 | U | 5.2 | 0.41 |
| 108-20-3 | Diisopropyl ether | 0.61 | U | 10 | 0.61 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.2 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.44 | U | 5.2 | 0.44 |
| 75-09-2 | Methylene Chloride | 0.76 | U | 5.2 | 0.76 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.43 | U | 10 | 0.43 |
| 91-20-3 | Naphthalene | 0.36 | U | 5.2 | 0.36 |
| 104-51-8 | n-Butylbenzene | 0.61 | U | 5.2 | 0.61 |
| 103-65-1 | N-Propylbenzene | 0.62 | U | 5.2 | 0.62 |
| 95-47-6 | o-Xylene | 0.59 | U | 5.2 | 0.59 |
| 135-98-8 | sec-Butylbenzene | 0.94 | U | 5.2 | 0.94 |
| 100-42-5 | Styrene | 0.41 | U | 5.2 | 0.41 |
| 994-05-8 | Tert-amyl methyl ether | 0.44 | U | 5.2 | 0.44 |
| 637-92-3 | Ethyl t-butyl ether | 0.42 | U | 5.2 | 0.42 |
| 127-18-4 | Tetrachloroethene | 0.83 | U | 5.2 | 0.83 |
| 108-88-3 | Toluene | 0.28 | U | 5.2 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.33 | U | 5.2 | 0.33 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.30 | U | 5.2 | 0.30 |
| 79-01-6 | Trichloroethene | 0.39 | U | 5.2 | 0.39 |
| 75-69-4 | Trichlorofluoromethane | 0.30 | U | 5.2 | 0.30 |
| 108-05-4 | Vinyl acetate | 0.68 | U | 10 | 0.68 |
| 75-01-4 | Vinyl chloride | 0.31 | U | 5.2 | 0.31 |
| 1330-20-7 | Xylenes, Total | 0.56 | U | 10 | 0.56 |
| 98-06-6 | tert-Butylbenzene | 0.69 | U | 5.2 | 0.69 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 146350.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 5.67(g) | Date Analyzed: 11/10/2015 19:44

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.55 | U | 5.2 | 0.55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 105 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 90 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 100 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 146351.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.55(g) | Date Analyzed: 11/10/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.53 | U | 5.3 | 0.53 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.68 | U | 5.3 | 0.68 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.33 | U | 5.3 | 0.33 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.44 | U | 5.3 | 0.44 |
| 75-34-3 | 1,1-Dichloroethane | 0.36 | U | 5.3 | 0.36 |
| 75-35-4 | 1,1-Dichloroethene | 0.84 | U | 5.3 | 0.84 |
| 563-58-6 | 1,1-Dichloropropene | 0.39 | U | 5.3 | 0.39 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.53 | U | 5.3 | 0.53 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.82 | U | 5.3 | 0.82 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.3 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.40 | U | 5.3 | 0.40 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.61 | U | 5.3 | 0.61 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.9 | U | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.36 | U | 5.3 | 0.36 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.3 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.3 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.73 | U | 5.3 | 0.73 |
| 142-28-9 | 1,3-Dichloropropane | 0.42 | U | 5.3 | 0.42 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.30 | U | 5.3 | 0.30 |
| 594-20-7 | 2,2-Dichloropropane | 0.46 | U | 5.3 | 0.46 |
| 78-93-3 | 2-Butanone (MEK) | 4.0 | J B | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.67 | U | 53 | 0.67 |
| 95-49-8 | 2-Chlorotoluene | 0.32 | U | 5.3 | 0.32 |
| 591-78-6 | 2-Hexanone | 0.72 | U | 21 | 0.72 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.3 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 5.3 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 6.0 | U | 21 | 6.0 |
| 71-43-2 | Benzene | 0.95 | U | 5.3 | 0.95 |
| 108-86-1 | Bromobenzene | 0.35 | U | 5.3 | 0.35 |
| 75-25-2 | Bromoform | 0.25 | U | 5.3 | 0.25 |
| 74-83-9 | Bromomethane | 0.44 | U | 5.3 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.62 | U | 5.3 | 0.62 |
| 56-23-5 | Carbon tetrachloride | 0.68 | U | 5.3 | 0.68 |
| 108-90-7 | Chlorobenzene | 0.56 | U | 5.3 | 0.56 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 146351.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.55(g) | Date Analyzed: 11/10/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.39 | U | 5.3 | 0.39 |
| 75-00-3 | Chloroethane | 0.46 | U | 5.3 | 0.46 |
| 67-66-3 | Chloroform | 0.38 | U | 5.3 | 0.38 |
| 74-87-3 | Chloromethane | 0.83 | U | 5.3 | 0.83 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.30 | U | 5.3 | 0.30 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.75 | U | 5.3 | 0.75 |
| 74-95-3 | Dibromomethane | 0.44 | U | 5.3 | 0.44 |
| 74-97-5 | Bromochloromethane | 0.35 | U | 5.3 | 0.35 |
| 75-71-8 | Dichlorodifluoromethane | 0.30 | U | 5.3 | 0.30 |
| 75-27-4 | Bromodichloromethane | 0.31 | U | 5.3 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.0 | U | 210 | 6.0 |
| 100-41-4 | Ethylbenzene | 0.34 | U | 5.3 | 0.34 |
| 87-68-3 | Hexachlorobutadiene | 0.42 | U | 5.3 | 0.42 |
| 108-20-3 | Diisopropyl ether | 0.62 | U | 11 | 0.62 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.3 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.45 | U | 5.3 | 0.45 |
| 75-09-2 | Methylene Chloride | 0.78 | U | 5.3 | 0.78 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.44 | U | 11 | 0.44 |
| 91-20-3 | Naphthalene | 0.37 | U | 5.3 | 0.37 |
| 104-51-8 | n-Butylbenzene | 0.62 | U | 5.3 | 0.62 |
| 103-65-1 | N-Propylbenzene | 0.63 | U | 5.3 | 0.63 |
| 95-47-6 | o-Xylene | 0.60 | U | 5.3 | 0.60 |
| 135-98-8 | sec-Butylbenzene | 0.96 | U | 5.3 | 0.96 |
| 100-42-5 | Styrene | 0.42 | U | 5.3 | 0.42 |
| 994-05-8 | Tert-amyl methyl ether | 0.45 | U | 5.3 | 0.45 |
| 637-92-3 | Ethyl t-butyl ether | 0.43 | U | 5.3 | 0.43 |
| 127-18-4 | Tetrachloroethene | 0.84 | U | 5.3 | 0.84 |
| 108-88-3 | Toluene | 0.29 | J | 5.3 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.34 | U | 5.3 | 0.34 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.31 | U | 5.3 | 0.31 |
| 79-01-6 | Trichloroethene | 0.40 | U | 5.3 | 0.40 |
| 75-69-4 | Trichlorofluoromethane | 0.31 | U | 5.3 | 0.31 |
| 108-05-4 | Vinyl acetate | 0.70 | U | 11 | 0.70 |
| 75-01-4 | Vinyl chloride | 0.32 | U | 5.3 | 0.32 |
| 1330-20-7 | Xylenes, Total | 0.57 | U | 11 | 0.57 |
| 98-06-6 | tert-Butylbenzene | 0.71 | U | 5.3 | 0.71 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 146351.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.55(g) | Date Analyzed: 11/10/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.56 | U | 5.3 | 0.56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 87 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 89 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 78 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 96 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 146365.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.55(g) | Date Analyzed: 11/11/2015 13:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.53 | U | 5.3 | 0.53 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.68 | U | 5.3 | 0.68 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.33 | U | 5.3 | 0.33 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.44 | U | 5.3 | 0.44 |
| 75-34-3 | 1,1-Dichloroethane | 0.36 | U | 5.3 | 0.36 |
| 75-35-4 | 1,1-Dichloroethene | 0.84 | U | 5.3 | 0.84 |
| 563-58-6 | 1,1-Dichloropropene | 0.39 | U | 5.3 | 0.39 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.53 | U | 5.3 | 0.53 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.82 | U | 5.3 | 0.82 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.3 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.40 | U | 5.3 | 0.40 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.61 | U | 5.3 | 0.61 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.9 | U | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.36 | U | 5.3 | 0.36 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.3 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.3 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.73 | U | 5.3 | 0.73 |
| 142-28-9 | 1,3-Dichloropropane | 0.42 | U | 5.3 | 0.42 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.29 | U | 5.3 | 0.29 |
| 594-20-7 | 2,2-Dichloropropane | 0.46 | U | 5.3 | 0.46 |
| 78-93-3 | 2-Butanone (MEK) | 3.6 | J B | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.67 | U | 53 | 0.67 |
| 95-49-8 | 2-Chlorotoluene | 0.32 | U | 5.3 | 0.32 |
| 591-78-6 | 2-Hexanone | 0.72 | U | 21 | 0.72 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.3 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 5.3 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 6.0 | U | 21 | 6.0 |
| 71-43-2 | Benzene | 0.95 | U | 5.3 | 0.95 |
| 108-86-1 | Bromobenzene | 0.35 | U | 5.3 | 0.35 |
| 75-25-2 | Bromoform | 0.25 | U | 5.3 | 0.25 |
| 74-83-9 | Bromomethane | 0.44 | U | 5.3 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.62 | U | 5.3 | 0.62 |
| 56-23-5 | Carbon tetrachloride | 0.68 | U | 5.3 | 0.68 |
| 108-90-7 | Chlorobenzene | 0.56 | U | 5.3 | 0.56 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 146365.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.55(g) | Date Analyzed: 11/11/2015 13:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.39 | U | 5.3 | 0.39 |
| 75-00-3 | Chloroethane | 0.46 | U | 5.3 | 0.46 |
| 67-66-3 | Chloroform | 0.38 | U | 5.3 | 0.38 |
| 74-87-3 | Chloromethane | 0.83 | U | 5.3 | 0.83 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.29 | U | 5.3 | 0.29 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.75 | U | 5.3 | 0.75 |
| 74-95-3 | Dibromomethane | 0.44 | U | 5.3 | 0.44 |
| 74-97-5 | Bromochloromethane | 0.35 | U | 5.3 | 0.35 |
| 75-71-8 | Dichlorodifluoromethane | 0.29 | U | 5.3 | 0.29 |
| 75-27-4 | Bromodichloromethane | 0.31 | U | 5.3 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.0 | U | 210 | 6.0 |
| 100-41-4 | Ethylbenzene | 0.34 | U | 5.3 | 0.34 |
| 87-68-3 | Hexachlorobutadiene | 0.42 | U | 5.3 | 0.42 |
| 108-20-3 | Diisopropyl ether | 0.62 | U | 11 | 0.62 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.3 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.45 | U | 5.3 | 0.45 |
| 75-09-2 | Methylene Chloride | 0.78 | U | 5.3 | 0.78 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.44 | U | 11 | 0.44 |
| 91-20-3 | Naphthalene | 0.37 | U | 5.3 | 0.37 |
| 104-51-8 | n-Butylbenzene | 0.62 | U | 5.3 | 0.62 |
| 103-65-1 | N-Propylbenzene | 0.63 | U | 5.3 | 0.63 |
| 95-47-6 | o-Xylene | 0.60 | U | 5.3 | 0.60 |
| 135-98-8 | sec-Butylbenzene | 0.96 | U | 5.3 | 0.96 |
| 100-42-5 | Styrene | 0.42 | U | 5.3 | 0.42 |
| 994-05-8 | Tert-amyl methyl ether | 0.45 | U | 5.3 | 0.45 |
| 637-92-3 | Ethyl t-butyl ether | 0.43 | U | 5.3 | 0.43 |
| 127-18-4 | Tetrachloroethene | 0.84 | U | 5.3 | 0.84 |
| 108-88-3 | Toluene | 0.28 | U | 5.3 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.34 | U | 5.3 | 0.34 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.31 | U | 5.3 | 0.31 |
| 79-01-6 | Trichloroethene | 0.40 | U | 5.3 | 0.40 |
| 75-69-4 | Trichlorofluoromethane | 0.31 | U | 5.3 | 0.31 |
| 108-05-4 | Vinyl acetate | 0.70 | U | 11 | 0.70 |
| 75-01-4 | Vinyl chloride | 0.32 | U | 5.3 | 0.32 |
| 1330-20-7 | Xylenes, Total | 0.57 | U | 11 | 0.57 |
| 98-06-6 | tert-Butylbenzene | 0.71 | U | 5.3 | 0.71 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 146365.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.55(g) | Date Analyzed: 11/11/2015 13:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.56 | U | 5.3 | 0.56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 102 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 89 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: TB-110515 | Lab Sample ID: 240-57597-16

Matrix: Water | Lab File ID: UXC4397.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.28 | U | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.44 | U | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.22 | U | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.30 | U | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 0.45 | U | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.37 | U | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.44 | U | 1.0 | 0.44 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.47 | U | 5.0 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.32 | U | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.41 | U | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.82 | U | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.25 | U | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 0.23 | U | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 0.25 | U | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.19 | U | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 0.19 | U | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.27 | U | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 0.26 | U | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 0.53 | U | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.59 | U | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 0.40 | U | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 0.48 | U | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 10 | 0.99 |
| 67-64-1 | Acetone | 0.94 | U | 10 | 0.94 |
| 71-43-2 | Benzene | 0.35 | U | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 0.35 | U | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 0.56 | U | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 0.44 | U | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.38 | U | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 0.43 | U | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 0.25 | U | 1.0 | 0.25 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: TB-110515 | Lab Sample ID: 240-57597-16

Matrix: Water | Lab File ID: UXC4397.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 1.0 | 0.43 |
| 75-00-3 | Chloroethane | 0.32 | U | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 0.25 | U | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 0.44 | U | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.26 | U | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.46 | U | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 0.42 | U | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.50 | U | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 4.9 | U | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 0.25 | U | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 0.35 | U | 1.0 | 0.35 |
| 108-20-3 | Diisopropyl ether | 0.50 | U | 10 | 0.50 |
| 98-82-8 | Isopropylbenzene | 0.35 | U | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 0.20 | U | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 0.33 | U | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.24 | U | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 0.45 | U | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 0.31 | U | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 0.40 | U | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 0.25 | U | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 0.48 | U | 1.0 | 0.48 |
| 100-42-5 | Styrene | 0.45 | U | 1.0 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.30 | U | 5.0 | 0.30 |
| 637-92-3 | Ethyl t-butyl ether | 0.23 | U | 5.0 | 0.23 |
| 127-18-4 | Tetrachloroethene | 0.31 | U | 1.0 | 0.31 |
| 108-88-3 | Toluene | 0.23 | U | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.30 | U | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.56 | U | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 0.22 | U | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 0.49 | U | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 0.41 | U | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 0.29 | U | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 0.52 | U | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 0.41 | U | 1.0 | 0.41 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: TB-110515 | Lab Sample ID: 240-57597-16

Matrix: Water | Lab File ID: UXC4397.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.32 | U | 1.0 | 0.32 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 85 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 97 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 90 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 99 | | 78-125 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50 Calibration End Date: 08/13/2015 20:42 Calibration ID: 30158

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-193150/20 | 144584.D |
| Level 2 | IC 240-193150/19 | 144583.D |
| Level 3 | IC 240-193150/18 | 144582.D |
| Level 4 | IC 240-193150/17 | 144581.D |
| Level 5 | IC 240-193150/16 | 144580.D |
| Level 6 | IC 240-193150/15 | 144579.D |
| Level 7 | IC 240-193150/14 | 144578.D |
| Level 8 | IC 240-193150/13 | 144577.D |
| Level 9 | IC 240-193150/12 | 144576.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Acetonitrile | 0.0270<br>0.0217 | 0.0199<br>0.0226 | 0.0223<br>0.0219 | 0.0211<br>0.0222 | 0.0231 | Ave | | 0.0224 | | | | 8.8 | | 15.0 | | | |
| 2-Chloro-1,3-butadiene | 0.3946<br>0.4330 | 0.4036<br>0.4520 | 0.4843<br>0.4367 | 0.4330<br>0.4384 | 0.4314 | Ave | | 0.4341 | | | | 6.0 | | 15.0 | | | |
| Diisopropyl ether | 0.1991<br>0.2087 | 0.1920<br>0.2170 | 0.2189<br>0.2067 | 0.2161<br>0.2066 | 0.2062 | Ave | | 0.2079 | | | | 4.2 | | 15.0 | | | |
| Ethyl t-butyl ether | 0.8061<br>0.8145 | 0.8124<br>0.8609 | 0.8459<br>0.8423 | 0.8104<br>0.8475 | 0.7961 | Ave | | 0.8262 | | | | 2.8 | | 15.0 | | | |
| Ethyl acetate | 0.2082<br>0.1555 | 0.1520<br>0.1600 | 0.1502<br>0.1572 | 0.1518<br>0.1578 | 0.1525 | Ave | | 0.1606 | | | | 11.3 | | 15.0 | | | |
| Propionitrile | 0.0291<br>0.0267 | 0.0294<br>0.0281 | 0.0277<br>0.0272 | 0.0279<br>0.0272 | 0.0264 | Ave | | 0.0278 | | | | 3.7 | | 15.0 | | | |
| Methacrylonitrile | 0.1071<br>0.1082 | 0.1056<br>0.1106 | 0.1086<br>0.1043 | 0.1077<br>0.1034 | 0.1075 | Ave | | 0.1070 | | | | 2.1 | | 15.0 | | | |
| Tert-amyl methyl ether | 0.6617<br>0.6962 | 0.6767<br>0.7331 | 0.7307<br>0.7077 | 0.7151<br>0.7140 | 0.6826 | Ave | | 0.7020 | | | | 3.5 | | 15.0 | | | |
| n-Butanol | 0.0138<br>0.0096 | 0.0095<br>0.0106 | 0.0097<br>0.0104 | 0.0114<br>0.0105 | 0.0089 | Ave | | 0.0105 | | | | 13.7 | | 15.0 | | | |
| Ethyl acrylate | 0.2316<br>0.2216 | 0.2219<br>0.2309 | 0.2178<br>0.2294 | 0.2151<br>0.2318 | 0.2122 | Ave | | 0.2236 | | | | 3.4 | | 15.0 | | | |
| Methyl methacrylate | 0.1524<br>0.1436 | 0.1363<br>0.1478 | 0.1383<br>0.1452 | 0.1404<br>0.1466 | 0.1384 | Ave | | 0.1432 | | | | 3.7 | | 15.0 | | | |
| 2-Nitropropane | +++++<br>0.0226 | 0.0209<br>0.0266 | 0.0198<br>0.0269 | 0.0195<br>0.0236 | 0.0194 | Ave | | 0.0224 | | | | 13.7 | | 15.0 | | | |
| n-Butyl acetate | +++++<br>0.3716 | +++++<br>0.3944 | 0.4011<br>0.3902 | 0.3921<br>0.3867 | 0.3596 | Ave | | 0.3851 | | | | 3.7 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50 Calibration End Date: 08/13/2015 20:42 Calibration ID: 30158

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1-Chlorohexane | 0.4929<br>0.4383 | 0.4320<br>0.4614 | 0.4729<br>0.4436 | 0.4373<br>0.4435 | 0.4315 | Ave | | 0.4504 | | | | 4.7 | | 15.0 | | | |
| Cyclohexanone | 0.0128<br>0.0108 | 0.0155<br>0.0119 | 0.0120<br>0.0114 | 0.0149<br>0.0114 | 0.0103 | Ave | | 0.0123 | | | | 14.4 | | 15.0 | | | |
| 1,2,3-Trimethylbenzene | 2.8189<br>3.0746 | 2.8992<br>3.2639 | 3.1500<br>3.1145 | 3.0040<br>3.0651 | 2.9168 | Ave | | 3.0341 | | | | 4.6 | | 15.0 | | | |
| 1,3,5-Trichlorobenzene | 1.1368<br>1.0908 | 1.0640<br>1.1383 | 1.1518<br>1.0486 | 1.1209<br>1.0391 | 1.0725 | Ave | | 1.0959 | | | | 3.9 | | 15.0 | | | |
| 2-Methylnaphthalene | 1.1258<br>1.2729 | 1.0980<br>1.4105 | 1.0984<br>1.2876 | 1.2519<br>1.2359 | 1.1966 | Ave | | 1.2197 | | | | 8.4 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50 Calibration End Date: 08/13/2015 20:42 Calibration ID: 30158

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-193150/20 | 144584.D |
| Level 2 | IC 240-193150/19 | 144583.D |
| Level 3 | IC 240-193150/18 | 144582.D |
| Level 4 | IC 240-193150/17 | 144581.D |
| Level 5 | IC 240-193150/16 | 144580.D |
| Level 6 | IC 240-193150/15 | 144579.D |
| Level 7 | IC 240-193150/14 | 144578.D |
| Level 8 | IC 240-193150/13 | 144577.D |
| Level 9 | IC 240-193150/12 | 144576.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Acetonitrile | FB | Ave | 7388<br>303462 | 10973<br>632599 | 30131<br>1256831 | 59051<br>1567842 | 126527 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| 2-Chloro-1,3-butadiene | FB | Ave | 10787<br>606721 | 22288<br>1265427 | 65473<br>2509599 | 121146<br>3096439 | 236791 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Diisopropyl ether | FB | Ave | 5444<br>292480 | 10602<br>607343 | 29596<br>1187824 | 60452<br>1459333 | 113175 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl t-butyl ether | FB | Ave | 22036<br>1141245 | 44869<br>2409991 | 114365<br>4840430 | 226764<br>5986416 | 436960 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl acetate | FB | Ave | 11385<br>435795 | 16786<br>895870 | 40612<br>1806750 | 84961<br>2228660 | 167431 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Propionitrile | FB | Ave | 7963<br>374531 | 16256<br>786705 | 37500<br>1563948 | 77984<br>1921197 | 145100 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| Methacrylonitrile | FB | Ave | 29273<br>1516021 | 58312<br>3095674 | 146830<br>5994798 | 301406<br>7302328 | 590265 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| Tert-amyl methyl ether | FB | Ave | 18089<br>975479 | 37373<br>2052126 | 98783<br>4066811 | 200083<br>5043426 | 374623 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| n-Butanol | CBZ | Ave | 5835<br>210549 | 8000<br>460443 | 20632<br>903775 | 49843<br>1135945 | 77408 | 25.0<br>1250 | 50.0<br>2500 | 125<br>5000 | 250<br>6250 | 500 |
| Ethyl acrylate | FB | Ave | 6332<br>310518 | 12255<br>646338 | 29451<br>1317957 | 60196<br>1637141 | 116465 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methyl methacrylate | FB | Ave | 8332<br>402313 | 15055<br>827639 | 37395<br>1668202 | 78576<br>2071659 | 151896 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| 2-Nitropropane | FB | Ave | +++++<br>63467 | 2313<br>149194 | 5344<br>309024 | 10920<br>333900 | 21282 | +++++<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| n-Butyl acetate | CBZ | Ave | +++++<br>325474 | +++++<br>683511 | 34073<br>1362203 | 68874<br>1673169 | 125276 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1-Chlorohexane | CBZ | Ave | 8362<br>383892 | 14627<br>799511 | 40171<br>1548771 | 76810<br>1919108 | 150333 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50 Calibration End Date: 08/13/2015 20:42 Calibration ID: 30158

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Cyclohexanone | CBZ | Ave | 2169<br>94475 | 5243<br>205876 | 10210<br>398420 | 26100<br>493296 | 35769 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| 1,2,3-Trimethylbenzene | DCB | Ave | 22714<br>1297881 | 46268<br>2748737 | 129836<br>5348625 | 256074<br>6584366 | 492514 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3,5-Trichlorobenzene | DCB | Ave | 9160<br>460463 | 16981<br>958626 | 47473<br>1800789 | 95549<br>2232228 | 181103 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Methylnaphthalene | DCB | Ave | 18143<br>1074702 | 35046<br>2375842 | 90547<br>4422384 | 213428<br>5309881 | 404111 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |

Curve Type Legend:
Ave = Average ISTD

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-196026/11 | 144943.D |
| Level 2 | IC 240-196026/10 | 144942.D |
| Level 3 | IC 240-196026/9 | 144941.D |
| Level 4 | IC 240-196026/8 | 144940.D |
| Level 5 | IC 240-196026/7 | 144939.D |
| Level 6 | ICIS 240-196026/6 | 144938.D |
| Level 7 | IC 240-196026/5 | 144937.D |
| Level 8 | IC 240-196026/4 | 144936.D |
| Level 9 | IC 240-196026/3 | 144935.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | +++++<br>0.2391 | +++++<br>0.2507 | 0.2470<br>0.2355 | 0.2306<br>0.2458 | 0.2468 | Ave | | 0.2422 | | | 0.1000 | 3.0 | | 15.0 | | | |
| Chloromethane | 0.3754<br>0.3823 | 0.4019<br>0.3850 | 0.3740<br>0.3707 | 0.3723<br>0.3714 | 0.3751 | Ave | | 0.3787 | | | 0.1000 | 2.6 | | 15.0 | | | |
| Vinyl chloride | 0.2535<br>0.2723 | 0.2872<br>0.2816 | 0.2779<br>0.2705 | 0.2715<br>0.2802 | 0.2773 | Ave | | 0.2747 | | | 0.1000 | 3.5 | | 15.0 | | | |
| Butadiene | 0.2421<br>0.3144 | 0.2898<br>0.3180 | 0.3006<br>0.3092 | 0.3016<br>0.3170 | 0.3102 | Ave | | 0.3003 | | | | 7.9 | | 15.0 | | | |
| Bromomethane | 0.1681<br>0.1234 | 0.1376<br>+++++ | 0.1308<br>+++++ | 0.1354<br>+++++ | 0.1348 | Ave | | 0.1383 | | | 0.0500 | 11.1 | | 15.0 | | | |
| Chloroethane | 0.1729<br>0.1392 | 0.1452<br>0.1270 | 0.1461<br>+++++ | 0.1422<br>+++++ | 0.1409 | Ave | | 0.1448 | | | | 9.6 | | 15.0 | | | |
| Dichlorofluoromethane | 0.3353<br>0.3214 | 0.3424<br>0.3238 | 0.3254<br>0.3013 | 0.3326<br>0.3083 | 0.3302 | Ave | | 0.3245 | | | | 4.0 | | 15.0 | | | |
| Trichlorofluoromethane | 0.2372<br>0.3085 | 0.2496<br>0.3130 | 0.2776<br>0.2931 | 0.2862<br>0.3038 | 0.2978 | Ave | | 0.2852 | | | 0.1000 | 9.2 | | 15.0 | | | |
| Ethyl ether | 0.2263<br>0.1633 | 0.1893<br>0.1738 | 0.1892<br>0.1620 | 0.1718<br>0.1640 | 0.1839 | Ave | | 0.1804 | | | | 11.2 | | 15.0 | | | |
| Acrolein | 0.0216<br>0.0171 | 0.0159<br>0.0175 | 0.0196<br>+++++ | 0.0166<br>+++++ | 0.0179 | Ave | | 0.0180 | | | | 10.9 | | 15.0 | | | |
| 1,1-Dichloroethene | 0.2079<br>0.2313 | 0.1781<br>0.2399 | 0.2268<br>0.2226 | 0.2292<br>0.2287 | 0.2305 | Ave | | 0.2217 | | | 0.1000 | 8.3 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.1337<br>0.1711 | 0.1398<br>0.1761 | 0.1720<br>0.1647 | 0.1739<br>0.1677 | 0.1631 | Ave | | 0.1625 | | | 0.0500 | 9.4 | | 15.0 | | | |
| Acetone | +++++<br>0.0712 | +++++<br>0.0711 | 0.1266<br>0.0622 | 0.0892<br>0.0631 | 0.0783 | Lin1 | 0.6353 | 0.0626 | | | 0.0500 | | | | 0.9980 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32  Calibration End Date: 09/02/2015 17:24  Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Iodomethane | 0.3031<br>0.3309 | 0.3161<br>0.3401 | 0.3156<br>0.3166 | 0.3107<br>0.3246 | 0.3168 | Ave | | 0.3194 | | | | 3.4 | | 15.0 | | | |
| Carbon disulfide | +++++<br>0.6938 | +++++<br>0.7470 | 0.5786<br>0.7111 | 0.6047<br>0.7368 | 0.6277 | Ave | | 0.6714 | | | 0.1000 | 10.0 | | 15.0 | | | |
| 3-Chloro-1-propene | +++++<br>0.1475 | +++++<br>0.1563 | 0.1275<br>0.1460 | 0.1372<br>0.1497 | 0.1368 | Ave | | 0.1430 | | | | 6.8 | | 15.0 | | | |
| Methyl acetate | 0.1615<br>0.1574 | 0.1594<br>0.1602 | 0.1513<br>0.1539 | 0.1584<br>0.1527 | 0.1554 | Ave | | 0.1567 | | | 0.1000 | 2.3 | | 15.0 | | | |
| Methylene Chloride | +++++<br>0.2588 | +++++<br>0.2535 | 0.5579<br>0.2285 | 0.3574<br>0.2329 | 0.3001 | Lin1 | 1.5559 | 0.2263 | | | 0.1000 | | | | 0.9990 | | 0.9900 |
| tert-Butyl alcohol (TBA) | 0.0223<br>0.0196 | 0.0192<br>0.0202 | 0.0186<br>0.0190 | 0.0196<br>0.0197 | 0.0203 | Ave | | 0.0198 | | | | 5.4 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.2358<br>0.2589 | 0.2329<br>0.2662 | 0.2511<br>0.2514 | 0.2599<br>0.2562 | 0.2510 | Ave | | 0.2515 | | | 0.1000 | 4.3 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.5407<br>0.5781 | 0.5287<br>0.5970 | 0.5676<br>0.5608 | 0.5729<br>0.5665 | 0.5729 | Ave | | 0.5650 | | | 0.1000 | 3.6 | | 15.0 | | | |
| Acrylonitrile | 0.0788<br>0.0761 | 0.0771<br>0.0770 | 0.0752<br>0.0728 | 0.0765<br>0.0724 | 0.0759 | Ave | | 0.0757 | | | | 2.7 | | 15.0 | | | |
| Hexane | +++++<br>0.0773 | +++++<br>0.0782 | 0.1040<br>0.0738 | 0.0811<br>0.0760 | 0.0806 | Ave | | 0.0816 | | | | 12.5 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.4672<br>0.5014 | 0.4509<br>0.5208 | 0.4998<br>0.4916 | 0.4999<br>0.5087 | 0.4920 | Ave | | 0.4925 | | | 0.1000 | 4.3 | | 15.0 | | | |
| Vinyl acetate | 0.0241<br>0.0308 | 0.0166<br>0.0307 | 0.0238<br>+++++ | 0.0268<br>+++++ | 0.0267 | Lin1 | -0.018 | 0.0303 | | | | | | | 0.9970 | | 0.9900 |
| 2,2-Dichloropropane | 0.1821<br>0.2315 | 0.1870<br>0.2370 | 0.2083<br>0.2070 | 0.2175<br>0.2121 | 0.2141 | Ave | | 0.2107 | | | | 8.5 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.2557<br>0.2665 | 0.2545<br>0.2811 | 0.2602<br>0.2612 | 0.2694<br>0.2684 | 0.2713 | Ave | | 0.2654 | | | 0.1000 | 3.2 | | 15.0 | | | |
| 2-Butanone (MEK) | +++++<br>0.0942 | 0.1118<br>0.0991 | 0.0952<br>0.0916 | 0.0894<br>0.0929 | 0.0959 | Ave | | 0.0962 | | | 0.0500 | 7.2 | | 15.0 | | | |
| Bromochloromethane | 0.0858<br>0.1003 | 0.0854<br>0.1062 | 0.0973<br>0.1000 | 0.0967<br>0.1004 | 0.0986 | Ave | | 0.0967 | | | | 7.1 | | 15.0 | | | |
| Tetrahydrofuran | +++++<br>0.0638 | 0.0727<br>0.0652 | 0.0687<br>0.0628 | 0.0716<br>0.0619 | 0.0652 | Ave | | 0.0665 | | | | 6.1 | | 15.0 | | | |
| Chloroform | 0.4109<br>0.3987 | 0.3844<br>0.4148 | 0.3916<br>0.3944 | 0.3857<br>0.4100 | 0.3904 | Ave | | 0.3979 | | | | 2.9 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.2815<br>0.3335 | 0.2455<br>0.3520 | 0.3014<br>0.3288 | 0.3164<br>0.3392 | 0.3164 | Ave | | 0.3127 | | | 0.1000 | 10.5 | | 15.0 | | | |
| Cyclohexane | 0.5387<br>0.6083 | 0.4961<br>0.6348 | 0.6113<br>0.5854 | 0.6257<br>0.6043 | 0.6056 | Ave | | 0.5900 | | | 0.1000 | 7.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32  Calibration End Date: 09/02/2015 17:24  Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon tetrachloride | 0.1974<br>0.2765 | 0.2070<br>0.2949 | 0.2344<br>0.2806 | 0.2508<br>0.2887 | 0.2562 | Ave | | 0.2541 | | | 0.1000 | 13.8 | | 15.0 | | | |
| 1,1-Dichloropropene | +++++<br>0.3352 | +++++<br>0.3502 | 0.3354<br>0.3258 | 0.3276<br>0.3378 | 0.3200 | Ave | | 0.3331 | | | | 3.0 | | 15.0 | | | |
| Benzene | 0.9461<br>1.0076 | 0.9369<br>1.0637 | 0.9831<br>0.9992 | 1.0297<br>1.0364 | 0.9873 | Ave | | 0.9989 | | | 0.5000 | 4.1 | | 15.0 | | | |
| Isobutyl alcohol | 0.0079<br>0.0070 | 0.0073<br>0.0070 | 0.0070<br>0.0065 | 0.0074<br>0.0064 | 0.0078 | Ave | | 0.0071 | | | | 7.3 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.2922<br>0.2806 | 0.2538<br>0.2902 | 0.2809<br>0.2746 | 0.2857<br>0.2775 | 0.2775 | Ave | | 0.2792 | | | 0.1000 | 4.0 | | 15.0 | | | |
| n-Heptane | 0.0495<br>0.0719 | 0.0549<br>0.0771 | 0.0671<br>0.0732 | 0.0733<br>0.0762 | 0.0704 | Ave | | 0.0682 | | | | 14.1 | | 15.0 | | | |
| Trichloroethene | 0.2322<br>0.2352 | 0.1925<br>0.2515 | 0.2312<br>0.2339 | 0.2474<br>0.2400 | 0.2383 | Ave | | 0.2336 | | | 0.2000 | 7.2 | | 15.0 | | | |
| Methylcyclohexane | 0.3977<br>0.5007 | 0.4180<br>0.5258 | 0.5131<br>0.4892 | 0.4996<br>0.5023 | 0.4970 | Ave | | 0.4826 | | | 0.1000 | 9.1 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.2351<br>0.2464 | 0.2304<br>0.2703 | 0.2458<br>0.2552 | 0.2519<br>0.2638 | 0.2445 | Ave | | 0.2493 | | | 0.1000 | 5.1 | | 15.0 | | | |
| Dibromomethane | 0.1138<br>0.1049 | 0.0928<br>0.1099 | 0.1030<br>0.1045 | 0.0989<br>0.1052 | 0.1030 | Ave | | 0.1040 | | | | 5.7 | | 15.0 | | | |
| 1,4-Dioxane | 0.0016<br>0.0015 | 0.0018<br>0.0016 | 0.0017<br>0.0016 | 0.0016<br>0.0016 | 0.0018 | Ave | | 0.0017 | | | | 5.4 | | 15.0 | | | |
| Bromodichloromethane | +++++<br>0.2583 | +++++<br>0.2789 | 0.2043<br>0.2740 | 0.2214<br>0.2839 | 0.2350 | Ave | | 0.2508 | | | 0.1000 | 12.3 | | 15.0 | | | |
| 2-Chloroethyl vinyl ether | 0.0880<br>0.0981 | 0.0738<br>0.1047 | 0.0849<br>0.1021 | 0.0908<br>0.1040 | 0.0933 | Ave | | 0.0933 | | | | 10.9 | | 15.0 | | | |
| cis-1,3-Dichloropropene | +++++<br>0.3137 | +++++<br>0.3452 | 0.2529<br>0.3385 | 0.2764<br>0.3487 | 0.2874 | Ave | | 0.3090 | | | 0.1500 | 12.2 | | 15.0 | | | |
| 4-Methyl-2-pentanone (MIBK) | 0.3475<br>0.3097 | 0.3025<br>0.3146 | 0.3102<br>0.2962 | 0.3082<br>0.2899 | 0.3133 | Ave | | 0.3102 | | | 0.0500 | 5.2 | | 15.0 | | | |
| Toluene | 1.6889<br>1.6108 | 1.5730<br>1.6794 | 1.5789<br>1.5858 | 1.6326<br>1.6292 | 1.5722 | Ave | | 1.6167 | | | 0.4000 | 2.8 | | 15.0 | | | |
| trans-1,3-Dichloropropene | +++++<br>0.3786 | +++++<br>0.4056 | 0.2981<br>0.3958 | 0.3438<br>0.4021 | 0.3521 | Ave | | 0.3680 | | | 0.1000 | 10.6 | | 15.0 | | | |
| Ethyl methacrylate | 0.3411<br>0.2949 | 0.2728<br>0.3134 | 0.2921<br>0.2986 | 0.3041<br>0.3000 | 0.2937 | Ave | | 0.3012 | | | | 6.1 | | 15.0 | | | |
| 1,1,2-Trichloroethane | 0.2387<br>0.2409 | 0.2020<br>0.2502 | 0.2341<br>0.2296 | 0.2342<br>0.2336 | 0.2367 | Ave | | 0.2333 | | | 0.1000 | 5.6 | | 15.0 | | | |
| Tetrachloroethene | 0.2604<br>0.2924 | 0.2503<br>0.3024 | 0.2921<br>0.2817 | 0.2907<br>0.2908 | 0.2885 | Ave | | 0.2833 | | | 0.2000 | 6.0 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32  Calibration End Date: 09/02/2015 17:24  Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,3-Dichloropropane | 0.4985<br>0.4202 | 0.3942<br>0.4386 | 0.3932<br>0.4193 | 0.4331<br>0.4249 | 0.4155 | Ave | | 0.4264 | | | | 7.3 | | 15.0 | | | |
| 2-Hexanone | 0.2259<br>0.1948 | 0.2056<br>0.1966 | 0.2059<br>0.1806 | 0.2139<br>0.1792 | 0.2041 | Ave | | 0.2007 | | | 0.0500 | 7.5 | | 15.0 | | | |
| Dibromochloromethane | +++++<br>0.2379 | 0.1676<br>0.2584 | 0.1731<br>0.2529 | 0.1970<br>0.2554 | 0.2047 | Lin1 | -0.324 | 0.2543 | | | | | | | 0.9990 | | 0.9900 |
| 1,2-Dibromoethane | 0.1915<br>0.2254 | 0.1731<br>0.2352 | 0.2140<br>0.2225 | 0.2199<br>0.2231 | 0.2195 | Ave | | 0.2138 | | | | 9.0 | | 15.0 | | | |
| Chlorobenzene | 0.8957<br>0.9330 | 0.9144<br>0.9681 | 0.9220<br>0.9247 | 0.9499<br>0.9450 | 0.9159 | Ave | | 0.9298 | | | 0.3000 | 2.3 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.2424<br>0.3183 | 0.2461<br>0.3292 | 0.2688<br>0.3070 | 0.2915<br>0.3071 | 0.3010 | Ave | | 0.2902 | | | | 10.7 | | 15.0 | | | |
| Ethylbenzene | 0.4967<br>0.5429 | 0.5117<br>0.5557 | 0.5343<br>0.5186 | 0.5496<br>0.5291 | 0.5275 | Ave | | 0.5296 | | | | 3.5 | | 15.0 | | | |
| m-Xylene & p-Xylene | 0.6493<br>0.6610 | 0.5935<br>0.6804 | 0.6630<br>0.6340 | 0.6529<br>0.6481 | 0.6392 | Ave | | 0.6468 | | | | 3.7 | | 15.0 | | | |
| o-Xylene | 0.6484<br>0.6613 | 0.5588<br>0.6783 | 0.6201<br>0.6144 | 0.6581<br>0.6145 | 0.6575 | Ave | | 0.6346 | | | | 5.7 | | 15.0 | | | |
| Styrene | 0.9630<br>0.9968 | 0.8800<br>1.0308 | 0.8828<br>0.9707 | 0.9401<br>0.9819 | 0.9540 | Ave | | 0.9556 | | | 0.3000 | 5.2 | | 15.0 | | | |
| Bromoform | 0.0617<br>0.1189 | 0.0736<br>0.1321 | 0.0749<br>0.1315 | 0.0937<br>0.1318 | 0.0977 | Lin1 | -0.136 | 0.1302 | | | 0.1000 | | | | 0.9970 | | 0.9900 |
| Isopropylbenzene | 1.6615<br>1.8186 | 1.5426<br>1.8815 | 1.7312<br>1.7266 | 1.7956<br>1.7460 | 1.7709 | Ave | | 1.7416 | | | 0.1000 | 5.6 | | 15.0 | | | |
| Bromobenzene | 0.7018<br>0.7269 | 0.6199<br>0.7522 | 0.6741<br>0.7263 | 0.7025<br>0.7372 | 0.6965 | Ave | | 0.7042 | | | | 5.6 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 0.5544<br>0.6344 | 0.5565<br>0.6397 | 0.5491<br>0.6163 | 0.5930<br>0.5999 | 0.6184 | Ave | | 0.5958 | | | 0.3000 | 5.9 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.1905<br>0.1847 | 0.1628<br>0.1813 | 0.1632<br>0.1748 | 0.1767<br>0.1723 | 0.1805 | Ave | | 0.1763 | | | | 5.3 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | 0.2675<br>0.2192 | 0.1583<br>0.2222 | 0.1934<br>0.2263 | 0.2038<br>0.2273 | 0.2186 | Ave | | 0.2152 | | | | 13.7 | | 15.0 | | | |
| N-Propylbenzene | 0.9746<br>1.0339 | 0.8541<br>1.0565 | 0.8793<br>1.0140 | 0.9898<br>1.0407 | 0.9489 | Ave | | 0.9769 | | | | 7.3 | | 15.0 | | | |
| 2-Chlorotoluene | 0.8414<br>0.8319 | 0.6742<br>0.8400 | 0.7804<br>0.8160 | 0.8063<br>0.8183 | 0.8038 | Ave | | 0.8014 | | | | 6.4 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 2.7121<br>3.3648 | 2.7168<br>3.3834 | 2.9008<br>3.2276 | 3.1601<br>3.2885 | 3.1601 | Ave | | 3.1016 | | | | 8.4 | | 15.0 | | | |
| 4-Chlorotoluene | 0.8033<br>0.7833 | 0.7373<br>0.8244 | 0.7337<br>0.7938 | 0.7735<br>0.8011 | 0.7475 | Ave | | 0.7776 | | | | 4.1 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| tert-Butylbenzene | 2.3374<br>2.9631 | 2.4133<br>3.0556 | 2.6550<br>2.8376 | 2.8314<br>2.8634 | 2.8090 | Ave | | 2.7518 | | | | 8.7 | | 15.0 | | | |
| 1,2,4-Trimethylbenzene | 2.7336<br>3.2868 | 2.6205<br>3.3941 | 2.8875<br>3.2352 | 3.0975<br>3.2908 | 3.1820 | Ave | | 3.0809 | | | | 8.8 | | 15.0 | | | |
| sec-Butylbenzene | 3.5842<br>4.5688 | 3.5902<br>4.6258 | 3.9612<br>4.3978 | 4.2686<br>4.4271 | 4.3250 | Ave | | 4.1943 | | | | 9.4 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.4872<br>1.4944 | 1.4279<br>1.5057 | 1.4500<br>1.4639 | 1.4918<br>1.4763 | 1.4693 | Ave | | 1.4741 | | | 0.6000 | 1.7 | | 15.0 | | | |
| p-Isopropyltoluene | 2.6700<br>3.5326 | 2.7312<br>3.5868 | 3.1681<br>3.3752 | 3.3032<br>3.4312 | 3.3784 | Ave | | 3.2418 | | | | 10.2 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 1.6796<br>1.4771 | 1.4525<br>1.4817 | 1.3997<br>1.4447 | 1.4789<br>1.4779 | 1.4661 | Ave | | 1.4842 | | | 0.5000 | 5.2 | | 15.0 | | | |
| n-Butylbenzene | 2.6703<br>3.3398 | 2.7947<br>3.4180 | 3.0393<br>3.3135 | 3.1433<br>3.3757 | 3.2352 | Ave | | 3.1478 | | | | 8.4 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.5577<br>1.4312 | 1.4056<br>1.4532 | 1.3574<br>1.3836 | 1.4146<br>1.3837 | 1.4022 | Ave | | 1.4210 | | | 0.4000 | 4.1 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | 0.0691<br>0.0964 | 0.0727<br>0.1007 | 0.0775<br>0.1005 | 0.0817<br>0.0987 | 0.0866 | Ave | | 0.0871 | | | 0.0500 | 14.3 | | 15.0 | | | |
| 1,2,4-Trichlorobenzene | 0.9903<br>0.9952 | 0.9066<br>0.9801 | 0.9495<br>0.9392 | 1.0303<br>0.9542 | 1.0144 | Ave | | 0.9733 | | | 0.2000 | 4.0 | | 15.0 | | | |
| Hexachlorobutadiene | 0.3960<br>0.5853 | 0.4364<br>0.5796 | 0.5790<br>0.5455 | 0.5879<br>0.5625 | 0.5837 | Ave | | 0.5396 | | | | 13.3 | | 15.0 | | | |
| Naphthalene | 2.5487<br>2.3198 | 2.2039<br>2.2590 | 2.1592<br>2.1437 | 2.2755<br>2.0812 | 2.3749 | Ave | | 2.2629 | | | | 6.2 | | 15.0 | | | |
| 1,2,3-Trichlorobenzene | 0.9121<br>0.9513 | 0.9280<br>0.9117 | 0.9440<br>0.8603 | 0.9768<br>0.8520 | 0.9646 | Ave | | 0.9223 | | | | 4.7 | | 15.0 | | | |
| Dibromofluoromethane (Surr) | +++++<br>0.2047 | +++++<br>0.2121 | 0.2454<br>0.1982 | 0.2289<br>0.2019 | 0.2161 | Ave | | 0.2154 | | | | 7.8 | | 15.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | +++++<br>0.2276 | +++++<br>0.2392 | 0.3027<br>0.2282 | 0.2781<br>0.2330 | 0.2521 | Ave | | 0.2516 | | | | 11.4 | | 15.0 | | | |
| Toluene-d8 (Surr) | +++++<br>1.4443 | +++++<br>1.4774 | 1.8704<br>1.3884 | 1.6510<br>1.4120 | 1.5274 | Ave | | 1.5387 | | | | 11.1 | | 15.0 | | | |
| 4-Bromofluorobenzene (Surr) | +++++<br>1.0832 | +++++<br>1.0727 | 1.4341<br>1.0549 | 1.2044<br>1.0662 | 1.1169 | Ave | | 1.1475 | | | | 11.9 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-196026/11 | 144943.D |
| Level 2 | IC 240-196026/10 | 144942.D |
| Level 3 | IC 240-196026/9 | 144941.D |
| Level 4 | IC 240-196026/8 | 144940.D |
| Level 5 | IC 240-196026/7 | 144939.D |
| Level 6 | ICIS 240-196026/6 | 144938.D |
| Level 7 | IC 240-196026/5 | 144937.D |
| Level 8 | IC 240-196026/4 | 144936.D |
| Level 9 | IC 240-196026/3 | 144935.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Dichlorodifluoromethane | FB | Ave | +++++<br>356137 | +++++<br>748933 | 36569<br>1466377 | 68169<br>1835598 | 147021 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Chloromethane | FB | Ave | 10383<br>569551 | 22625<br>1150096 | 55373<br>2307997 | 110043<br>2773342 | 223410 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Vinyl chloride | FB | Ave | 7011<br>405713 | 16172<br>841040 | 41149<br>1684326 | 80229<br>2092495 | 165146 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Butadiene | FB | Ave | 6696<br>468393 | 16314<br>949979 | 44512<br>1925445 | 89146<br>2367224 | 184770 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Bromomethane | FB | Ave | 4648<br>183791 | 7747<br>+++++ | 19368<br>+++++ | 40006<br>+++++ | 80268 | 1.00<br>50.0 | 2.00<br>+++++ | 5.00<br>+++++ | 10.0<br>+++++ | 20.0 |
| Chloroethane | FB | Ave | 4781<br>207308 | 8175<br>379413 | 21638<br>+++++ | 42019<br>+++++ | 83922 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>+++++ | 10.0<br>+++++ | 20.0 |
| Dichlorofluoromethane | FB | Ave | 9272<br>478742 | 19276<br>967265 | 48180<br>1875722 | 98295<br>2301848 | 196686 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Trichlorofluoromethane | FB | Ave | 6559<br>459614 | 14053<br>934958 | 41107<br>1824715 | 84588<br>2268583 | 177390 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl ether | FB | Ave | 6258<br>243336 | 10659<br>519027 | 28019<br>1008376 | 50769<br>1224410 | 109526 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Acrolein | FB | Ave | 2993<br>127696 | 4483<br>261405 | 14479<br>+++++ | 24467<br>+++++ | 53257 | 5.00<br>250 | 10.0<br>500 | 25.0<br>+++++ | 50.0<br>+++++ | 100 |
| 1,1-Dichloroethene | FB | Ave | 5750<br>344590 | 10027<br>716734 | 33584<br>1386108 | 67737<br>1707262 | 137258 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroetha ne | FB | Ave | 3699<br>254932 | 7868<br>526084 | 25468<br>1025667 | 51393<br>1251857 | 97164 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Acetone | FB | Lin1 | +++++<br>212159 | +++++<br>424676 | 37501<br>774275 | 52703<br>941871 | 93287 | +++++<br>100 | +++++<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Iodomethane | FB | Ave | 8383<br>493006 | 17794<br>1015933 | 46725<br>1971089 | 91832<br>2423809 | 188704 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Carbon disulfide | FB | Ave | +++++<br>1033515 | +++++<br>2231244 | 85670<br>4427381 | 178714<br>5501396 | 373862 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 3-Chloro-1-propene | FB | Ave | +++++<br>219767 | +++++<br>467005 | 18885<br>908999 | 40539<br>1117889 | 81474 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methyl acetate | FB | Ave | 22330<br>1172266 | 44885<br>2392801 | 112036<br>4789789 | 234059<br>5701867 | 462647 | 5.00<br>250 | 10.0<br>500 | 25.0<br>1000 | 50.0<br>1250 | 100 |
| Methylene Chloride | FB | Lin1 | +++++<br>385484 | +++++<br>757088 | 82611<br>1422440 | 105644<br>1738961 | 178739 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| tert-Butyl alcohol (TBA) | FB | Ave | 6178<br>292151 | 10809<br>601929 | 27588<br>1184581 | 57843<br>1471581 | 120716 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| trans-1,2-Dichloroethene | FB | Ave | 6522<br>385670 | 13114<br>795293 | 37178<br>1565230 | 76822<br>1912734 | 149499 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methyl tert-butyl ether | FB | Ave | 14953<br>861134 | 29764<br>1783409 | 84049<br>3491615 | 169317<br>4230179 | 341246 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Acrylonitrile | FB | Ave | 21785<br>1133959 | 43399<br>2301252 | 111279<br>4534284 | 225966<br>5402864 | 451855 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| Hexane | FB | Ave | +++++<br>115126 | +++++<br>233508 | 15392<br>459224 | 23963<br>567539 | 48009 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1-Dichloroethane | FB | Ave | 12922<br>746990 | 25387<br>1555732 | 74007<br>3060770 | 147740<br>3798439 | 293058 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Vinyl acetate | FB | Lin1 | 667<br>45823 | 932<br>91562 | 3519<br>+++++ | 7932<br>+++++ | 15884 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>+++++ | 10.0<br>+++++ | 20.0 |
| 2,2-Dichloropropane | FB | Ave | 5036<br>344804 | 10529<br>708031 | 30835<br>1289042 | 64289<br>1583856 | 127532 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| cis-1,2-Dichloroethene | FB | Ave | 7073<br>396968 | 14329<br>839540 | 38524<br>1626156 | 79621<br>2004348 | 161592 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Butanone (MEK) | FB | Ave | +++++<br>280552 | 12585<br>591753 | 28186<br>1140829 | 52850<br>1386909 | 114195 | +++++<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Bromochloromethane | FB | Ave | 2372<br>149423 | 4806<br>317360 | 14400<br>622349 | 28593<br>749740 | 58707 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Tetrahydrofuran | FB | Ave | +++++<br>189961 | 8190<br>389802 | 20359<br>781838 | 42328<br>924687 | 77626 | +++++<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Chloroform | FB | Ave | 11364<br>593941 | 21643<br>1239162 | 57984<br>2455954 | 113989<br>3061654 | 232521 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,1-Trichloroethane | FB | Ave | 7784<br>496863 | 13820<br>1051452 | 44626<br>2047378 | 93504<br>2532683 | 188459 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Cyclohexane | FB | Ave | 14900<br>906203 | 27932<br>1896235 | 90514<br>3644821 | 184933<br>4512267 | 360725 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Carbon tetrachloride | FB | Ave | 5460<br>411902 | 11655<br>880737 | 34704<br>1747228 | 74120<br>2155848 | 152621 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1-Dichloropropene | FB | Ave | +++++<br>499401 | +++++<br>1046178 | 49664<br>2028361 | 96828<br>2521880 | 190594 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Benzene | FB | Ave | 26167<br>1500996 | 52745<br>3177421 | 145567<br>6221614 | 304344<br>7738061 | 588026 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Isobutyl alcohol | FB | Ave | 5460<br>261074 | 10213<br>525940 | 25933<br>1010906 | 54805<br>1189785 | 116205 | 25.0<br>1250 | 50.0<br>2500 | 125<br>5000 | 250<br>6250 | 500 |
| 1,2-Dichloroethane | FB | Ave | 8081<br>418078 | 14287<br>866965 | 41595<br>1709845 | 84448<br>2072165 | 165251 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| n-Heptane | FB | Ave | 1370<br>107045 | 3089<br>230164 | 9941<br>455483 | 21652<br>568917 | 41942 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Trichloroethene | FB | Ave | 6423<br>350358 | 10835<br>751379 | 34227<br>1456606 | 73115<br>1792315 | 141942 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methylcyclohexane | FB | Ave | 10999<br>745915 | 23534<br>1570602 | 75969<br>3045944 | 147660<br>3750421 | 296027 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dichloropropane | FB | Ave | 6501<br>367015 | 12972<br>807304 | 36398<br>1589013 | 74445<br>1969959 | 145632 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Dibromomethane | FB | Ave | 3146<br>156268 | 5223<br>328160 | 15254<br>650540 | 29240<br>785166 | 61356 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,4-Dioxane | FB | Ave | 896<br>45240 | 2050<br>96254 | 4894<br>200368 | 9657<br>245495 | 21042 | 20.0<br>1000 | 40.0<br>2000 | 100<br>4000 | 200<br>5000 | 400 |
| Bromodichloromethane | FB | Ave | +++++<br>384799 | +++++<br>832998 | 30251<br>1706175 | 65422<br>2119797 | 139991 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Chloroethyl vinyl ether | FB | Ave | 4867<br>292355 | 8305<br>625520 | 25131<br>1271843 | 53690<br>1553516 | 111110 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| cis-1,3-Dichloropropene | FB | Ave | +++++<br>467262 | +++++<br>1031197 | 37447<br>2107636 | 81678<br>2603965 | 171161 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | CBZ | Ave | 11429<br>563617 | 20282<br>1178361 | 55788<br>2332635 | 109762<br>2766526 | 227626 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Toluene | CBZ | Ave | 27774<br>1465877 | 52739<br>3144782 | 141953<br>6243364 | 290701<br>7772646 | 571155 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| trans-1,3-Dichloropropene | CBZ | Ave | +++++<br>344567 | +++++<br>759613 | 26806<br>1558216 | 61223<br>1918393 | 127919 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl methacrylate | CBZ | Ave | 5609<br>268334 | 9146<br>586878 | 26258<br>1175568 | 54145<br>1431045 | 106690 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,2-Trichloroethane | CBZ | Ave | 3926<br>219261 | 6772<br>468456 | 21044<br>904034 | 41703<br>1114274 | 86000 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Tetrachloroethene | CBZ | Ave | 4283<br>266087 | 8391<br>566338 | 26258<br>1109089 | 51758<br>1387568 | 104805 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3-Dichloropropane | CBZ | Ave | 8198<br>382372 | 13218<br>821259 | 35354<br>1650749 | 77111<br>2027259 | 150930 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Hexanone | CBZ | Ave | 7429<br>354469 | 13785<br>736413 | 37025<br>1421673 | 76163<br>1709783 | 148306 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Dibromochloromethane | CBZ | Lin1 | +++++<br>216458 | 5618<br>483817 | 15565<br>995515 | 35077<br>1218638 | 74370 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1,2-Dibromoethane | CBZ | Ave | 3150<br>205128 | 5804<br>440477 | 19238<br>875919 | 39164<br>1064183 | 79732 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Chlorobenzene | CBZ | Ave | 14730<br>849090 | 30658<br>1812785 | 82892<br>3640404 | 169143<br>4508577 | 332714 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,1,2-Tetrachloroethane | CBZ | Ave | 3986<br>289664 | 8252<br>616483 | 24169<br>1208539 | 51910<br>1465339 | 109363 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethylbenzene | CBZ | Ave | 8169<br>494070 | 17157<br>1040646 | 48037<br>2041599 | 97870<br>2524079 | 191624 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| m-Xylene & p-Xylene | CBZ | Ave | 10677<br>601578 | 19898<br>1274050 | 59612<br>2495890 | 116259<br>3091955 | 232208 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| o-Xylene | CBZ | Ave | 10663<br>601783 | 18737<br>1270093 | 55753<br>2418978 | 117189<br>2931823 | 238863 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Styrene | CBZ | Ave | 15837<br>907146 | 29506<br>1930215 | 79372<br>3821764 | 167399<br>4684740 | 346563 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Bromoform | CBZ | Lin1 | 1015<br>108179 | 2467<br>247433 | 6730<br>517661 | 16693<br>628700 | 35497 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Isopropylbenzene | CBZ | Ave | 27324<br>1654997 | 51721<br>3523233 | 155645<br>6797585 | 319732<br>8330012 | 643348 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Bromobenzene | DCB | Ave | 5573<br>295347 | 9956<br>626961 | 28765<br>1217640 | 58476<br>1492683 | 116607 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,2,2-Tetrachloroethane | DCB | Ave | 4403<br>257748 | 8937<br>533250 | 23432<br>1033232 | 49363<br>1214765 | 103530 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,3-Trichloropropane | DCB | Ave | 1513<br>75031 | 2614<br>151158 | 6965<br>293104 | 14705<br>348980 | 30218 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| trans-1,4-Dichloro-2-butene | DCB | Ave | 2124<br>89070 | 2543<br>185204 | 8252<br>379325 | 16966<br>460351 | 36604 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| N-Propylbenzene | DCB | Ave | 7740<br>420052 | 13717<br>880640 | 37524<br>1699968 | 82391<br>2107190 | 158855 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Chlorotoluene | DCB | Ave | 6682<br>337984 | 10827<br>700147 | 33303<br>1368070 | 67113<br>1656963 | 134563 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3,5-Trimethylbenzene | DCB | Ave | 21538<br>1367130 | 43631<br>2820139 | 123786<br>5410976 | 263032<br>6658839 | 529049 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 4-Chlorotoluene | DCB | Ave | 6379<br>318271 | 11841<br>687196 | 31307<br>1330849 | 64384<br>1622168 | 125148 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| tert-Butylbenzene | DCB | Ave | 18562<br>1203890 | 38757<br>2546917 | 113296<br>4757117 | 235674<br>5798117 | 470266 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,4-Trimethylbenzene | DCB | Ave | 21708<br>1335433 | 42084<br>2829044 | 123218<br>5423776 | 257826<br>6663540 | 532704 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| sec-Butylbenzene | DCB | Ave | 28463<br>1856287 | 57658<br>3855727 | 169034<br>7372707 | 355305<br>8964364 | 724063 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3-Dichlorobenzene | DCB | Ave | 11810<br>607174 | 22932<br>1255024 | 61875<br>2454144 | 124176<br>2989387 | 245980 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| p-Isopropyltoluene | DCB | Ave | 21203<br>1435272 | 43863<br>2989705 | 135190<br>5658412 | 274948<br>6947777 | 565586 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,4-Dichlorobenzene | DCB | Ave | 13338<br>600156 | 23326<br>1235039 | 59729<br>2421962 | 123097<br>2992653 | 245441 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| n-Butylbenzene | DCB | Ave | 21206<br>1356939 | 44882<br>2849004 | 129694<br>5555001 | 261638<br>6835319 | 541621 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dichlorobenzene | DCB | Ave | 12370<br>581486 | 22574<br>1211298 | 57923<br>2319608 | 117745<br>2801754 | 234745 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | DCB | Ave | 549<br>39179 | 1167<br>83939 | 3309<br>168459 | 6800<br>199946 | 14490 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,4-Trichlorobenzene | DCB | Ave | 7864<br>404356 | 14560<br>816947 | 40518<br>1574505 | 85761<br>1932210 | 169826 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Hexachlorobutadiene | DCB | Ave | 3145<br>237801 | 7009<br>483093 | 24708<br>914525 | 48937<br>1138913 | 97726 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Naphthalene | DCB | Ave | 20240<br>942540 | 35394<br>1882955 | 92141<br>3593926 | 189406<br>4214177 | 397593 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,3-Trichlorobenzene | DCB | Ave | 7243<br>386525 | 14903<br>759901 | 40282<br>1442298 | 81306<br>1725103 | 161494 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Dibromofluoromethane (Surr) | FB | Ave | +++++<br>304994 | +++++<br>633562 | 36336<br>1234341 | 67662<br>1507714 | 128718 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Ave | +++++<br>339100 | +++++<br>714630 | 44817<br>1421037 | 82196<br>1739346 | 150166 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Toluene-d8 (Surr) | CBZ | Ave | +++++<br>1314337 | +++++<br>2766497 | 168163<br>5466311 | 293978<br>6736325 | 554888 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 4-Bromofluorobenzene (Surr) | DCB | Ave | +++++<br>440117 | +++++<br>894108 | 61195<br>1768529 | 100246<br>2158955 | 186984 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin1 = Linear 1/conc ISTD |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-204067/7 | UXC3829A.D |
| Level 2 | STD8260 240-204067/6 | UXC3828A.D |
| Level 3 | STD8260 240-204067/5 | UXC3827A.D |
| Level 4 | STD8260 240-204067/4 | UXC3826A.D |
| Level 5 | STD8260 240-204067/3 | UXC3825A.D |
| Level 6 | STD8260 240-204067/2 | UXC3824A.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | 0.3763 0.4630 | 0.3685 | 0.4347 | 0.4095 | 0.4415 | Ave | | 0.4156 | | | 0.1000 | 9.1 | | 15.0 | | | |
| Chloromethane | 0.3134 0.3006 | 0.2822 | 0.3097 | 0.2949 | 0.3057 | Ave | | 0.3011 | | | 0.1000 | 3.8 | | 15.0 | | | |
| Vinyl chloride | 0.3653 0.3315 | 0.3066 | 0.3371 | 0.3133 | 0.3248 | Ave | | 0.3298 | | | 0.1000 | 6.3 | | 15.0 | | | |
| Butadiene | 0.3149 0.2921 | 0.2799 | 0.2958 | 0.2812 | 0.2944 | Ave | | 0.2931 | | | | 4.3 | | 15.0 | | | |
| Bromomethane | 0.1768 0.1785 | 0.1856 | 0.1888 | 0.1677 | 0.1766 | Ave | | 0.1790 | | | 0.0500 | 4.2 | | 15.0 | | | |
| Chloroethane | 0.2170 0.1827 | 0.1810 | 0.1946 | 0.1792 | 0.1864 | Ave | | 0.1902 | | | | 7.5 | | 15.0 | | | |
| Dichlorofluoromethane | 0.5491 0.5336 | 0.5008 | 0.5423 | 0.5081 | 0.5307 | Ave | | 0.5274 | | | | 3.6 | | 15.0 | | | |
| Trichlorofluoromethane | 0.4290 0.5156 | 0.4293 | 0.4745 | 0.4580 | 0.4761 | Ave | | 0.4637 | | | 0.1000 | 7.1 | | 15.0 | | | |
| Ethyl ether | 0.2524 0.2058 | 0.2022 | 0.2174 | 0.2184 | 0.2166 | Ave | | 0.2188 | | | | 8.1 | | 15.0 | | | |
| Acrolein | 0.0194 0.0208 | 0.0195 | 0.0213 | 0.0216 | 0.0207 | Ave | | 0.0206 | | | | 4.5 | | 15.0 | | | |
| 1,1-Dichloroethene | 0.3192 0.3103 | 0.3026 | 0.3175 | 0.3087 | 0.3189 | Ave | | 0.3129 | | | 0.1000 | 2.2 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.2468 0.2655 | 0.2417 | 0.2611 | 0.2571 | 0.2662 | Ave | | 0.2564 | | | 0.0500 | 3.9 | | 15.0 | | | |
| Acetone | 0.1494 0.0756 | 0.1046 | 0.0806 | 0.0779 | 0.0760 | Lin1 | 0.1360 | 0.0728 | | | 0.0500 | | | | 0.9990 | | 0.9900 |
| Iodomethane | 0.4808 0.4866 | 0.4808 | 0.4981 | 0.4871 | 0.4948 | Ave | | 0.4880 | | | | 1.5 | | 15.0 | | | |
| Carbon disulfide | 0.8042 0.9355 | 0.7777 | 0.8325 | 0.8658 | 0.9300 | Ave | | 0.8576 | | | 0.1000 | 7.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 3-Chloro-1-propene | 0.1574<br>0.1826 | 0.1531 | 0.1725 | 0.1821 | 0.1872 | Ave | | 0.1725 | | | | 8.3 | | 15.0 | | | |
| Methyl acetate | 0.1529<br>0.1529 | 0.1449 | 0.1499 | 0.1485 | 0.1560 | Ave | | 0.1509 | | | 0.1000 | 2.6 | | 15.0 | | | |
| Methylene Chloride | 0.6600<br>0.3263 | 0.4591 | 0.4025 | 0.3526 | 0.3465 | Lin1 | 0.3272 | 0.3223 | | | 0.1000 | | | | 0.9990 | | 0.9900 |
| tert-Butyl alcohol (TBA) | 0.0126<br>0.0126 | 0.0137 | 0.0130 | 0.0136 | 0.0145 | Ave | | 0.0133 | | | | 5.5 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.7083<br>0.8027 | 0.7349 | 0.7554 | 0.7782 | 0.8151 | Ave | | 0.7657 | | | 0.1000 | 5.3 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.3546<br>0.3362 | 0.3350 | 0.3511 | 0.3434 | 0.3491 | Ave | | 0.3449 | | | 0.1000 | 2.3 | | 15.0 | | | |
| Acrylonitrile | 0.0807<br>0.0839 | 0.0777 | 0.0809 | 0.0821 | 0.0849 | Ave | | 0.0817 | | | | 3.1 | | 15.0 | | | |
| Hexane | 0.0897<br>0.0895 | 0.0775 | 0.0959 | 0.0883 | 0.0933 | Ave | | 0.0890 | | | | 7.1 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.6004<br>0.5990 | 0.5656 | 0.6055 | 0.5988 | 0.6163 | Ave | | 0.5976 | | | 0.1000 | 2.9 | | 15.0 | | | |
| Vinyl acetate | 0.2891<br>0.3663 | 0.2933 | 0.3167 | 0.3246 | 0.3559 | Ave | | 0.3243 | | | | 9.8 | | 15.0 | | | |
| 2,2-Dichloropropane | 0.2966<br>0.3299 | 0.3036 | 0.3180 | 0.3379 | 0.3501 | Ave | | 0.3227 | | | | 6.4 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.3967<br>0.3605 | 0.3546 | 0.3747 | 0.3635 | 0.3760 | Ave | | 0.3710 | | | 0.1000 | 4.1 | | 15.0 | | | |
| 2-Butanone (MEK) | 0.0890<br>0.0927 | 0.0889 | 0.0877 | 0.0928 | 0.0965 | Ave | | 0.0913 | | | 0.0500 | 3.7 | | 15.0 | | | |
| Bromochloromethane | 0.1697<br>0.1681 | 0.1526 | 0.1693 | 0.1646 | 0.1743 | Ave | | 0.1664 | | | | 4.5 | | 15.0 | | | |
| Tetrahydrofuran | 0.0551<br>0.0522 | 0.0497 | 0.0500 | 0.0514 | 0.0520 | Ave | | 0.0518 | | | | 3.7 | | 15.0 | | | |
| Chloroform | 0.6342<br>0.6461 | 0.6330 | 0.6483 | 0.6338 | 0.6635 | Ave | | 0.6432 | | | | 1.9 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.4416<br>0.5089 | 0.4542 | 0.4774 | 0.5040 | 0.5219 | Ave | | 0.4847 | | | 0.1000 | 6.6 | | 15.0 | | | |
| Cyclohexane | 0.5510<br>0.5788 | 0.5254 | 0.5645 | 0.5688 | 0.5952 | Ave | | 0.5639 | | | 0.1000 | 4.3 | | 15.0 | | | |
| Carbon tetrachloride | 0.3590<br>0.4615 | 0.3535 | 0.4034 | 0.4258 | 0.4608 | Ave | | 0.4107 | | | 0.1000 | 11.6 | | 15.0 | | | |
| 1,1-Dichloropropene | 0.4475<br>0.4891 | 0.4505 | 0.4742 | 0.4878 | 0.4988 | Ave | | 0.4747 | | | | 4.5 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Isobutyl alcohol | 0.0045<br>0.0052 | 0.0043 | 0.0050 | 0.0051 | 0.0054 | Ave | | 0.0049 | | | | 8.3 | | 15.0 | | | |
| Benzene | 1.4285<br>1.4410 | 1.3752 | 1.3957 | 1.3924 | 1.4478 | Ave | | 1.4134 | | | 0.5000 | 2.1 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.5642<br>0.5505 | 0.5164 | 0.5367 | 0.5384 | 0.5567 | Ave | | 0.5438 | | | 0.1000 | 3.1 | | 15.0 | | | |
| n-Heptane | 0.0744<br>0.0875 | 0.0820 | 0.0810 | 0.0857 | 0.0897 | Ave | | 0.0834 | | | | 6.6 | | 15.0 | | | |
| Trichloroethene | 0.3624<br>0.3674 | 0.3397 | 0.3792 | 0.3617 | 0.3763 | Ave | | 0.3644 | | | 0.2000 | 3.9 | | 15.0 | | | |
| Methylcyclohexane | 0.5343<br>0.5582 | 0.5050 | 0.5318 | 0.5510 | 0.5699 | Ave | | 0.5417 | | | 0.1000 | 4.3 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.2971<br>0.3003 | 0.2855 | 0.3021 | 0.2932 | 0.3103 | Ave | | 0.2981 | | | 0.1000 | 2.8 | | 15.0 | | | |
| Dibromomethane | 0.1762<br>0.1826 | 0.1628 | 0.1757 | 0.1777 | 0.1863 | Ave | | 0.1769 | | | | 4.5 | | 15.0 | | | |
| 1,4-Dioxane | 0.0027<br>0.0022 | 0.0028 | 0.0024 | 0.0026 | 0.0028 | Ave | | 0.0026 | | | | 9.4 | | 15.0 | | | |
| Bromodichloromethane | 0.3493<br>0.4531 | 0.3322 | 0.3819 | 0.4022 | 0.4403 | Ave | | 0.3932 | | | 0.1000 | 12.3 | | 15.0 | | | |
| 2-Chloroethyl vinyl ether | 0.1365<br>0.1714 | 0.1411 | 0.1536 | 0.1565 | 0.1661 | Ave | | 0.1542 | | | | 8.8 | | 15.0 | | | |
| cis-1,3-Dichloropropene | 0.3364<br>0.4977 | 0.3567 | 0.4034 | 0.4277 | 0.4767 | Lin1 | -0.234 | 0.4896 | | | 0.1500 | | | | 0.9970 | | 0.9900 |
| 4-Methyl-2-pentanone (MIBK) | 0.1433<br>0.1852 | 0.1446 | 0.1677 | 0.1669 | 0.1802 | Ave | | 0.1647 | | | 0.0500 | 10.6 | | 15.0 | | | |
| Toluene | 1.9429<br>2.0302 | 1.8468 | 1.9605 | 1.9532 | 2.0824 | Ave | | 1.9693 | | | 0.4000 | 4.1 | | 15.0 | | | |
| trans-1,3-Dichloropropene | 0.3301<br>0.5505 | 0.3500 | 0.4184 | 0.4518 | 0.5264 | Lin1 | -0.328 | 0.5405 | | | 0.1000 | | | | 0.9960 | | 0.9900 |
| Ethyl methacrylate | 0.2971<br>0.4364 | 0.3159 | 0.3628 | 0.3705 | 0.4242 | Lin1 | -0.201 | 0.4307 | | | | | | | 0.9970 | | 0.9900 |
| 1,1,2-Trichloroethane | 0.3319<br>0.3389 | 0.3188 | 0.3304 | 0.3242 | 0.3461 | Ave | | 0.3317 | | | 0.1000 | 3.0 | | 15.0 | | | |
| Tetrachloroethene | 0.3571<br>0.3590 | 0.3460 | 0.3546 | 0.3549 | 0.3728 | Ave | | 0.3574 | | | 0.2000 | 2.5 | | 15.0 | | | |
| 1,3-Dichloropropane | 0.6595<br>0.6405 | 0.6138 | 0.6291 | 0.6236 | 0.6430 | Ave | | 0.6349 | | | | 2.5 | | 15.0 | | | |
| 2-Hexanone | 0.1368<br>0.1734 | 0.1326 | 0.1431 | 0.1652 | 0.1705 | Ave | | 0.1536 | | | 0.0500 | 11.8 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dibromochloromethane | 0.2548<br>0.3871 | 0.2497 | 0.2924 | 0.3190 | 0.3650 | Lin1 | -0.213 | 0.3778 | | | | | | | 0.9950 | | 0.9900 |
| 1,2-Dibromoethane | 0.3173<br>0.3432 | 0.2888 | 0.3287 | 0.3283 | 0.3440 | Ave | | 0.3251 | | | | 6.3 | | 15.0 | | | |
| Chlorobenzene | 1.3001<br>1.3022 | 1.2131 | 1.2671 | 1.2315 | 1.2845 | Ave | | 1.2664 | | | 0.3000 | 2.9 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.3166<br>0.4285 | 0.3268 | 0.3688 | 0.3781 | 0.4230 | Ave | | 0.3736 | | | | 12.5 | | 15.0 | | | |
| Ethylbenzene | 0.6139<br>0.6752 | 0.6272 | 0.6531 | 0.6630 | 0.6854 | Ave | | 0.6530 | | | | 4.2 | | 15.0 | | | |
| m-Xylene & p-Xylene | 1.6840<br>1.8801 | 1.6715 | 1.7805 | 1.7879 | 1.9286 | Ave | | 1.7888 | | | | 5.7 | | 15.0 | | | |
| o-Xylene | 0.7577<br>0.8160 | 0.7146 | 0.8018 | 0.7672 | 0.8323 | Ave | | 0.7816 | | | | 5.6 | | 15.0 | | | |
| Styrene | 1.2019<br>1.5179 | 1.2259 | 1.3763 | 1.3604 | 1.4960 | Ave | | 1.3631 | | | 0.3000 | 9.7 | | 15.0 | | | |
| Bromoform | 0.1273<br>0.1961 | 0.1074 | 0.1343 | 0.1474 | 0.1793 | Lin1 | -0.127 | 0.1882 | | | 0.1000 | | | | 0.9900 | | 0.9900 |
| Isopropylbenzene | 1.9913<br>2.0593 | 1.9528 | 2.1800 | 2.1276 | 2.3334 | Ave | | 2.1074 | | | 0.1000 | 6.6 | | 15.0 | | | |
| Bromobenzene | 0.9480<br>0.9172 | 0.8989 | 0.9318 | 0.9579 | 0.9573 | Ave | | 0.9352 | | | | 2.5 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 0.6367<br>0.6917 | 0.6612 | 0.6532 | 0.6916 | 0.6953 | Ave | | 0.6716 | | | 0.3000 | 3.7 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.2311<br>0.2455 | 0.2233 | 0.2502 | 0.2524 | 0.2593 | Ave | | 0.2436 | | | | 5.6 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | 0.1847<br>0.2273 | 0.1712 | 0.1873 | 0.2097 | 0.2190 | Ave | | 0.1999 | | | | 11.0 | | 15.0 | | | |
| N-Propylbenzene | 0.9712<br>1.0332 | 0.9264 | 1.0226 | 1.0490 | 1.0539 | Ave | | 1.0094 | | | | 5.0 | | 15.0 | | | |
| 2-Chlorotoluene | 0.8756<br>0.8951 | 0.8629 | 0.9068 | 0.9045 | 0.9220 | Ave | | 0.8945 | | | | 2.4 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 2.9650<br>3.6195 | 3.0996 | 3.3535 | 3.4351 | 3.5925 | Ave | | 3.3442 | | | | 7.9 | | 15.0 | | | |
| 4-Chlorotoluene | 3.4227<br>3.5911 | 3.2938 | 3.4666 | 3.4283 | 3.5772 | Ave | | 3.4633 | | | | 3.2 | | 15.0 | | | |
| tert-Butylbenzene | 2.6379<br>3.0255 | 2.5257 | 2.7206 | 2.8657 | 2.9264 | Ave | | 2.7836 | | | | 6.8 | | 15.0 | | | |
| 1,2,4-Trimethylbenzene | 3.2473<br>3.3932 | 3.1596 | 3.5475 | 3.6375 | 3.7459 | Ave | | 3.4552 | | | | 6.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| sec-Butylbenzene | 3.5593<br>3.6545 | 3.5760 | 3.8326 | 3.9468 | 4.1043 | Ave | | 3.7789 | | | | 5.8 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.9172<br>1.8494 | 1.7772 | 1.8392 | 1.8390 | 1.8648 | Ave | | 1.8478 | | | 0.6000 | 2.4 | | 15.0 | | | |
| p-Isopropyltoluene | 3.1956<br>3.3525 | 3.2534 | 3.4587 | 3.6196 | 3.7475 | Ave | | 3.4379 | | | | 6.2 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 2.0031<br>1.8854 | 1.9004 | 1.8823 | 1.8816 | 1.9099 | Ave | | 1.9104 | | | 0.5000 | 2.4 | | 15.0 | | | |
| n-Butylbenzene | 2.6219<br>3.0543 | 2.5547 | 2.7935 | 2.9524 | 3.0366 | Ave | | 2.8356 | | | | 7.5 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.8119<br>1.7466 | 1.7372 | 1.7233 | 1.7882 | 1.7748 | Ave | | 1.7637 | | | 0.4000 | 1.9 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | 0.0810<br>0.1175 | 0.0782 | 0.0887 | 0.1005 | 0.1127 | Lin1 | -0.061 | 0.1154 | | | 0.0500 | | | | 0.9960 | | 0.9900 |
| 1,2,4-Trichlorobenzene | 1.2025<br>1.1327 | 1.0994 | 1.1644 | 1.1753 | 1.1939 | Ave | | 1.1614 | | | 0.2000 | 3.4 | | 15.0 | | | |
| Hexachlorobutadiene | 0.5980<br>0.4967 | 0.5848 | 0.5440 | 0.5646 | 0.5659 | Ave | | 0.5590 | | | | 6.4 | | 15.0 | | | |
| Naphthalene | 2.1201<br>2.7163 | 2.2108 | 2.4032 | 2.6263 | 2.8052 | Ave | | 2.4803 | | | | 11.3 | | 15.0 | | | |
| 1,2,3-Trichlorobenzene | 1.1193<br>0.9977 | 0.9897 | 1.0555 | 1.0704 | 1.0863 | Ave | | 1.0532 | | | | 4.8 | | 15.0 | | | |
| Dibromofluoromethane (Surr) | 0.3005<br>0.3229 | 0.3208 | 0.3302 | 0.3349 | 0.3342 | Ave | | 0.3239 | | | | 4.0 | | 15.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | 0.4565<br>0.4542 | 0.4715 | 0.4518 | 0.4649 | 0.4642 | Ave | | 0.4605 | | | | 1.6 | | 15.0 | | | |
| Toluene-d8 (Surr) | 1.7070<br>1.8254 | 1.7324 | 1.7232 | 1.7818 | 1.8396 | Ave | | 1.7682 | | | | 3.2 | | 15.0 | | | |
| 4-Bromofluorobenzene (Surr) | 0.6725<br>0.6778 | 0.6451 | 0.6646 | 0.6672 | 0.6928 | Ave | | 0.6700 | | | | 2.4 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-204067/7 | UXC3829A.D |
| Level 2 | STD8260 240-204067/6 | UXC3828A.D |
| Level 3 | STD8260 240-204067/5 | UXC3827A.D |
| Level 4 | STD8260 240-204067/4 | UXC3826A.D |
| Level 5 | STD8260 240-204067/3 | UXC3825A.D |
| Level 6 | STD8260 240-204067/2 | UXC3824A.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Dichlorodifluoromethane | FB | Ave | 81451<br>4169025 | 163271 | 484098 | 926577 | 1995610 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Chloromethane | FB | Ave | 67835<br>2706728 | 125034 | 344799 | 667309 | 1381689 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Vinyl chloride | FB | Ave | 79073<br>2984679 | 135866 | 375367 | 708978 | 1467924 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Butadiene | FB | Ave | 68162<br>2630204 | 124013 | 329378 | 636245 | 1330914 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Bromomethane | FB | Ave | 38268<br>1607387 | 82232 | 210253 | 379551 | 798140 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Chloroethane | FB | Ave | 46977<br>1645135 | 80183 | 216709 | 405428 | 842721 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Dichlorofluoromethane | FB | Ave | 118849<br>4804244 | 221934 | 603891 | 1149672 | 2398842 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Trichlorofluoromethane | FB | Ave | 92850<br>4642420 | 190211 | 528323 | 1036260 | 2152000 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl ether | FB | Ave | 54631<br>1853290 | 89588 | 242132 | 494203 | 979228 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Acrolein | FB | Ave | 20991<br>937291 | 43213 | 118532 | 244644 | 468940 | 5.00<br>200 | 10.0 | 25.0 | 50.0 | 100 |
| 1,1-Dichloroethene | FB | Ave | 69093<br>2793947 | 134101 | 353528 | 698533 | 1441441 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroetha ne | FB | Ave | 53411<br>2390251 | 107089 | 290718 | 581817 | 1203179 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Acetone | FB | Lin1 | 64664<br>1361377 | 92744 | 179545 | 352749 | 687363 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Iodomethane | FB | Ave | 104063<br>4381841 | 213034 | 554660 | 1102088 | 2236420 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Carbon disulfide | FB | Ave | 174064<br>8423139 | 344599 | 927015 | 1959159 | 4203773 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 3-Chloro-1-propene | FB | Ave | 34075<br>1644594 | 67837 | 192042 | 412012 | 845942 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Methyl acetate | FB | Ave | 165477<br>6883090 | 321065 | 834592 | 1680575 | 3524921 | 5.00<br>200 | 10.0 | 25.0 | 50.0 | 100 |
| Methylene Chloride | FB | Lin1 | 142848<br>2937905 | 203456 | 448202 | 797761 | 1566439 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| tert-Butyl alcohol (TBA) | FB | Ave | 27273<br>1137904 | 60601 | 144403 | 307461 | 654244 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Methyl tert-butyl ether | FB | Ave | 153316<br>7227654 | 325626 | 841123 | 1760760 | 3684219 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| trans-1,2-Dichloroethene | FB | Ave | 76749<br>3026934 | 148462 | 390993 | 776906 | 1577937 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Acrylonitrile | FB | Ave | 174707<br>7558362 | 344389 | 901060 | 1857836 | 3837278 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Hexane | FB | Ave | 19413<br>805934 | 34338 | 106736 | 199837 | 421670 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1-Dichloroethane | FB | Ave | 129958<br>5393295 | 250613 | 674191 | 1354846 | 2785967 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Vinyl acetate | FB | Ave | 62581<br>3298089 | 129955 | 352626 | 734562 | 1608862 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2,2-Dichloropropane | FB | Ave | 64191<br>2970157 | 134515 | 354139 | 764552 | 1582302 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| cis-1,2-Dichloroethene | FB | Ave | 85871<br>3246257 | 157145 | 417246 | 822590 | 1699426 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Butanone (MEK) | FB | Ave | 38548<br>1668902 | 78751 | 195279 | 420122 | 872711 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Bromochloromethane | FB | Ave | 36730<br>1513516 | 67603 | 188519 | 372494 | 787744 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Tetrahydrofuran | FB | Ave | 23863<br>939205 | 44077 | 111449 | 232646 | 470504 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Chloroform | FB | Ave | 137284<br>5817632 | 280504 | 721918 | 1434085 | 2999162 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1,1-Trichloroethane | FB | Ave | 95594<br>4582510 | 201278 | 531601 | 1140336 | 2358943 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Cyclohexane | FB | Ave | 119256<br>5211339 | 232804 | 628590 | 1287004 | 2690588 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Carbon tetrachloride | FB | Ave | 77705<br>4155489 | 156632 | 449183 | 963488 | 2082898 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1-Dichloropropene | FB | Ave | 96854<br>4404131 | 199635 | 528051 | 1103828 | 2254841 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Isobutyl alcohol | FB | Ave | 24610<br>1174973 | 48123 | 140453 | 285810 | 612767 | 25.0<br>1000 | 50.0 | 125 | 250 | 500 |
| Benzene | FB | Ave | 309192<br>12974974 | 609356 | 1554073 | 3150718 | 6544023 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| 1,2-Dichloroethane | FB | Ave | 122119<br>4956680 | 228806 | 597601 | 1218319 | 2516349 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| n-Heptane | FB | Ave | 16106<br>787951 | 36356 | 90211 | 194014 | 405660 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Trichloroethene | FB | Ave | 78449<br>3308170 | 150514 | 422217 | 818405 | 1700845 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Methylcyclohexane | FB | Ave | 115643<br>5025981 | 223765 | 592115 | 1246817 | 2576081 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dichloropropane | FB | Ave | 64306<br>2703745 | 126512 | 336384 | 663367 | 1402405 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Dibromomethane | FB | Ave | 38148<br>1644188 | 72140 | 195647 | 401976 | 842177 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,4-Dioxane | FB | Ave | 11745<br>394929 | 24678 | 54334 | 115663 | 256040 | 20.0<br>800 | 40.0 | 100 | 200 | 400 |
| Bromodichloromethane | FB | Ave | 75615<br>4079952 | 147196 | 425207 | 910094 | 1990381 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Chloroethyl vinyl ether | FB | Ave | 59095<br>3087358 | 125047 | 341986 | 708402 | 1501781 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| cis-1,3-Dichloropropene | FB | Lin1 | 72806<br>4480949 | 158047 | 449240 | 967784 | 2154909 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | FB | Ave | 62056<br>3334443 | 128150 | 373529 | 755456 | 1629069 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Toluene | CBZ | Ave | 314750<br>13742655 | 613284 | 1662261 | 3337602 | 7008232 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| trans-1,3-Dichloropropene | CBZ | Lin1 | 53472<br>3726284 | 116215 | 354711 | 772009 | 1771713 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl methacrylate | CBZ | Lin1 | 48136<br>2953995 | 104902 | 307595 | 633162 | 1427604 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1,2-Trichloroethane | CBZ | Ave | 53764<br>2294126 | 105866 | 280175 | 553963 | 1164673 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Tetrachloroethene | CBZ | Ave | 57848<br>2430253 | 114893 | 300665 | 606536 | 1254556 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,3-Dichloropropane | CBZ | Ave | 106836<br>4335594 | 203818 | 533411 | 1065638 | 2163937 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Hexanone | CBZ | Ave | 44317<br>2347499 | 88069 | 242700 | 564530 | 1147465 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Dibromochloromethane | CBZ | Lin1 | 41280<br>2620558 | 82932 | 247907 | 545050 | 1228320 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dibromoethane | CBZ | Ave | 51396<br>2323252 | 95909 | 278692 | 561082 | 1157734 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Chlorobenzene | CBZ | Ave | 210609<br>8814971 | 402833 | 1074366 | 2104395 | 4322692 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| 1,1,1,2-Tetrachloroethane | CBZ | Ave | 51295<br>2900830 | 108508 | 312729 | 646033 | 1423538 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethylbenzene | CBZ | Ave | 99450<br>4570595 | 208277 | 553744 | 1132933 | 2306500 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| m-Xylene & p-Xylene | CBZ | Ave | 272799<br>12726899 | 555073 | 1509645 | 3055080 | 6490588 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| o-Xylene | CBZ | Ave | 122753<br>5523858 | 237315 | 679855 | 1311023 | 2801087 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Styrene | CBZ | Ave | 194713<br>10275242 | 407108 | 1166925 | 2324660 | 5034612 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Bromoform | CBZ | Lin1 | 20620<br>1327764 | 35682 | 113841 | 251934 | 603362 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Isopropylbenzene | CBZ | Ave | 322582<br>13939847 | 648484 | 1848388 | 3635569 | 7852789 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Bromobenzene | DCB | Ave | 84164<br>3502266 | 162589 | 434694 | 862350 | 1781313 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1,2,2-Tetrachloroethane | DCB | Ave | 56530<br>2641114 | 119597 | 304704 | 622565 | 1293771 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2,3-Trichloropropane | DCB | Ave | 20513<br>937347 | 40388 | 116709 | 227197 | 482566 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| trans-1,4-Dichloro-2-butene | DCB | Ave | 16396<br>868034 | 30971 | 87358 | 188753 | 407436 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| N-Propylbenzene | DCB | Ave | 86221<br>3945407 | 167571 | 477015 | 944360 | 1961128 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Chlorotoluene | DCB | Ave | 77733<br>3417835 | 156077 | 422997 | 814284 | 1715734 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,3,5-Trimethylbenzene | DCB | Ave | 263228<br>13821025 | 560662 | 1564358 | 3092405 | 6685007 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 4-Chlorotoluene | DCB | Ave | 303864<br>13712263 | 595781 | 1617111 | 3086235 | 6656394 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| tert-Butylbenzene | DCB | Ave | 234189<br>11552554 | 456850 | 1269117 | 2579794 | 5445541 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2,4-Trimethylbenzene | DCB | Ave | 288296<br>12956940 | 571516 | 1654850 | 3274598 | 6970355 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| sec-Butylbenzene | DCB | Ave | 315988<br>13954530 | 646817 | 1787847 | 3553029 | 7637226 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,3-Dichlorobenzene | DCB | Ave | 170206<br>7061962 | 321467 | 857954 | 1655516 | 3470085 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| p-Isopropyltoluene | DCB | Ave | 283699<br>12801430 | 588470 | 1613422 | 3258495 | 6973409 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,4-Dichlorobenzene | DCB | Ave | 177831<br>7199208 | 343752 | 878063 | 1693832 | 3554000 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| n-Butylbenzene | DCB | Ave | 232769<br>11662713 | 462098 | 1303137 | 2657792 | 5650481 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dichlorobenzene | DCB | Ave | 160858<br>6669488 | 314219 | 803878 | 1609759 | 3302623 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | DCB | Lin1 | 7188<br>448494 | 14136 | 41393 | 90507 | 209792 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2,4-Trichlorobenzene | DCB | Ave | 106759<br>4325166 | 198864 | 543198 | 1057994 | 2221645 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Hexachlorobutadiene | DCB | Ave | 53090<br>1896613 | 105782 | 253747 | 508257 | 1053091 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Naphthalene | DCB | Ave | 188220<br>10371948 | 399889 | 1121039 | 2364237 | 5220004 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2,3-Trichlorobenzene | DCB | Ave | 99374<br>3809727 | 179009 | 492383 | 963589 | 2021371 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Dibromofluoromethane (Surr) | FB | Ave | 65035<br>2907632 | 142131 | 367643 | 757894 | 1510819 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Ave | 98807<br>4089270 | 208921 | 503129 | 1051942 | 2098332 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Toluene-d8 (Surr) | CBZ | Ave | 276528<br>12356342 | 575295 | 1461082 | 3044709 | 6191118 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 4-Bromofluorobenzene (Surr) | CBZ | Ave | 108950<br>4587838 | 214241 | 563481 | 1140121 | 2331520 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin1 = Linear 1/conc ISTD |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 15:44 Calibration End Date: 10/28/2015 17:37 Calibration ID: 31161

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD 240-204067/13 | UXC3836.D |
| Level 2 | STD 240-204067/12 | UXC3835.D |
| Level 3 | STD 240-204067/11 | UXC3834.D |
| Level 4 | STD 240-204067/10 | UXC3833.D |
| Level 5 | STD 240-204067/9 | UXC3832.D |
| Level 6 | STD 240-204067/8 | UXC3831.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Acetonitrile | 0.0252 0.0219 | 0.0212 | 0.0233 | 0.0227 | 0.0226 | Ave | | 0.0228 | | | | 6.0 | | 15.0 | | | |
| Diisopropyl ether | 0.2683 0.2949 | 0.2607 | 0.2786 | 0.2899 | 0.2896 | Ave | | 0.2803 | | | | 4.8 | | 15.0 | | | |
| 2-Chloro-1,3-butadiene | 0.5072 0.5586 | 0.4581 | 0.5204 | 0.5317 | 0.5354 | Ave | | 0.5186 | | | | 6.6 | | 15.0 | | | |
| Ethyl t-butyl ether | 0.7639 0.8817 | 0.7574 | 0.8038 | 0.8303 | 0.8343 | Ave | | 0.8119 | | | | 5.8 | | 15.0 | | | |
| Ethyl acetate | 0.1723 0.1688 | 0.1526 | 0.1659 | 0.1628 | 0.1640 | Ave | | 0.1644 | | | | 4.1 | | 15.0 | | | |
| Propionitrile | 0.0300 0.0295 | 0.0271 | 0.0298 | 0.0298 | 0.0295 | Ave | | 0.0293 | | | | 3.7 | | 15.0 | | | |
| Methacrylonitrile | 0.1341 0.1392 | 0.1230 | 0.1302 | 0.1352 | 0.1327 | Ave | | 0.1324 | | | | 4.2 | | 15.0 | | | |
| Tert-amyl methyl ether | 0.6054 0.7084 | 0.5887 | 0.6367 | 0.6663 | 0.6719 | Ave | | 0.6462 | | | | 6.9 | | 15.0 | | | |
| n-Butanol | 0.0033 0.0038 | 0.0033 | 0.0034 | 0.0036 | 0.0038 | Ave | | 0.0035 | | | | 6.7 | | 15.0 | | | |
| Ethyl acrylate | 0.2445 0.2717 | 0.2373 | 0.2318 | 0.2579 | 0.2659 | Ave | | 0.2515 | | | | 6.4 | | 15.0 | | | |
| Methyl methacrylate | 0.1740 0.2003 | 0.1580 | 0.1764 | 0.1847 | 0.1885 | Ave | | 0.1803 | | | | 8.0 | | 15.0 | | | |
| 2-Nitropropane | 0.0283 0.0326 | 0.0236 | 0.0254 | 0.0265 | 0.0282 | Ave | | 0.0274 | | | | 11.3 | | 15.0 | | | |
| n-Butyl acetate | 0.2122 0.2685 | 0.2003 | 0.2057 | 0.2348 | 0.2498 | Ave | | 0.2286 | | | | 11.8 | | 15.0 | | | |
| 1-Chlorohexane | 0.4981 0.5513 | 0.4815 | 0.5004 | 0.5112 | 0.5405 | Ave | | 0.5138 | | | | 5.2 | | 15.0 | | | |
| Cyclohexanone | 0.0101 0.0133 | 0.0128 | 0.0128 | 0.0134 | 0.0144 | Ave | | 0.0128 | | | | 11.4 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 15:44 Calibration End Date: 10/28/2015 17:37 Calibration ID: 31161

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,2,3-Trimethylbenzene | 3.2126<br>3.7399 | 3.1433 | 3.4898 | 3.6396 | 3.7890 | Ave | | 3.5024 | | | | 7.8 | | 15.0 | | | |
| 1,3,5-Trichlorobenzene | 1.2561<br>1.2723 | 1.2439 | 1.2909 | 1.2690 | 1.3139 | Ave | | 1.2743 | | | | 2.0 | | 15.0 | | | |
| 2-Methylnaphthalene | 1.1675<br>1.4101 | 1.2466 | 1.4182 | 1.5971 | 1.7053 | Ave | | 1.4241 | | | | 14.3 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 15:44 Calibration End Date: 10/28/2015 17:37 Calibration ID: 31161

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD 240-204067/13 | UXC3836.D |
| Level 2 | STD 240-204067/12 | UXC3835.D |
| Level 3 | STD 240-204067/11 | UXC3834.D |
| Level 4 | STD 240-204067/10 | UXC3833.D |
| Level 5 | STD 240-204067/9 | UXC3832.D |
| Level 6 | STD 240-204067/8 | UXC3831.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Acetonitrile | FB | Ave | 55244<br>1935513 | 95625 | 258385 | 503778 | 1008896 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Diisopropyl ether | FB | Ave | 58924<br>2605872 | 117750 | 309424 | 642956 | 1295070 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Chloro-1,3-butadiene | FB | Ave | 111396<br>4935815 | 206907 | 578033 | 1179324 | 2394276 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl t-butyl ether | FB | Ave | 167760<br>7790456 | 342093 | 892694 | 1841635 | 3730781 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl acetate | FB | Ave | 75682<br>2982411 | 137897 | 368473 | 722018 | 1466626 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Propionitrile | FB | Ave | 65895<br>2606576 | 122305 | 330798 | 661402 | 1319574 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Methacrylonitrile | FB | Ave | 294563<br>12299460 | 555372 | 1445910 | 2998475 | 5932635 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Tert-amyl methyl ether | FB | Ave | 132956<br>6259304 | 265917 | 707158 | 1477825 | 3004628 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| n-Butanol | FB | Ave | 17977<br>832780 | 37465 | 93374 | 197738 | 426385 | 25.0<br>1000 | 50.0 | 125 | 250 | 500 |
| Ethyl acrylate | FB | Ave | 53688<br>2400896 | 107197 | 257438 | 572006 | 1188866 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Methyl methacrylate | FB | Ave | 76439<br>3539016 | 142738 | 391859 | 819344 | 1686044 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| 2-Nitropropane | FB | Ave | 12430<br>576148 | 21341 | 56329 | 117671 | 252142 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| n-Butyl acetate | FB | Ave | 46612<br>2371970 | 90494 | 228480 | 520730 | 1117091 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1-Chlorohexane | CBZ | Ave | 81498<br>3785206 | 158853 | 425093 | 889870 | 1841015 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Cyclohexanone | DCB | Ave | 8827<br>475638 | 21525 | 55001 | 119809 | 250164 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| 1,2,3-Trimethylbenzene | DCB | Ave | 280496<br>13327169 | 529556 | 1504300 | 3243980 | 6573591 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 15:44 Calibration End Date: 10/28/2015 17:37 Calibration ID: 31161

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| 1,3,5-Trichlorobenzene | DCB | Ave | 109673<br>4533718 | 209556 | 556440 | 1131040 | 2279554 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Methylnaphthalene | DCB | Ave | 203871<br>10050047 | 420018 | 1222651 | 2846954 | 5916979 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |

Curve Type Legend:
Ave = Average ISTD

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-196026/13 Calibration Date: 09/02/2015 18:07

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 144945.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2422 | 0.2902 | 0.1000 | 0.0240 | 0.0200 | 19.8 | 20.0 |
| Chloromethane | Ave | 0.3787 | 0.4264 | 0.1000 | 0.0225 | 0.0200 | 12.6 | 20.0 |
| Vinyl chloride | Ave | 0.2747 | 0.2991 | 0.1000 | 0.0218 | 0.0200 | 8.9 | 20.0 |
| Butadiene | Ave | 0.3003 | 0.3280 | | 0.0218 | 0.0200 | 9.2 | 20.0 |
| Bromomethane | Ave | 0.1383 | 0.1509 | 0.0500 | 0.0218 | 0.0200 | 9.1 | 20.0 |
| Chloroethane | Ave | 0.1448 | 0.1548 | 0.0500 | 0.0214 | 0.0200 | 6.9 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3245 | 0.3543 | | 0.0218 | 0.0200 | 9.2 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2852 | 0.3259 | 0.1000 | 0.0229 | 0.0200 | 14.3 | 20.0 |
| Ethyl ether | Ave | 0.1804 | 0.2164 | | 0.0240 | 0.0200 | 19.9 | 20.0 |
| Acrolein | Ave | 0.0180 | 0.0235 | | 0.130 | 0.100 | 30.4* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2217 | 0.2411 | 0.1000 | 0.0218 | 0.0200 | 8.8 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1625 | 0.1744 | 0.0500 | 0.0215 | 0.0200 | 7.4 | 20.0 |
| Acetone | Lin1 | | 0.0867 | 0.0500 | 0.0453 | 0.0400 | 13.2 | 20.0 |
| Iodomethane | Ave | 0.3194 | 0.3613 | | 0.0226 | 0.0200 | 13.1 | 20.0 |
| Carbon disulfide | Ave | 0.6714 | 0.6482 | 0.1000 | 0.0193 | 0.0200 | -3.4 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1430 | 0.1394 | | 0.0195 | 0.0200 | -2.5 | 20.0 |
| Methyl acetate | Ave | 0.1567 | 0.1835 | 0.1000 | 0.117 | 0.100 | 17.1 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3235 | 0.1000 | 0.0217 | 0.0200 | 8.6 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0198 | 0.0221 | | 0.223 | 0.200 | 11.3 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2515 | 0.2789 | 0.1000 | 0.0222 | 0.0200 | 10.9 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.5650 | 0.6276 | 0.1000 | 0.0222 | 0.0200 | 11.1 | 20.0 |
| Acrylonitrile | Ave | 0.0757 | 0.0862 | | 0.228 | 0.200 | 13.8 | 20.0 |
| Hexane | Ave | 0.0816 | 0.0814 | | 0.0200 | 0.0200 | -0.1 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4925 | 0.5567 | 0.1000 | 0.0226 | 0.0200 | 13.0 | 20.0 |
| Vinyl acetate | Lin1 | | 0.0217 | | 0.0149 | 0.0200 | -25.6* | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2107 | 0.2260 | | 0.0214 | 0.0200 | 7.2 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2654 | 0.2924 | 0.1000 | 0.0220 | 0.0200 | 10.2 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0962 | 0.1043 | 0.0500 | 0.0433 | 0.0400 | 8.4 | 20.0 |
| Bromochloromethane | Ave | 0.0967 | 0.1063 | | 0.0220 | 0.0200 | 9.9 | 20.0 |
| Tetrahydrofuran | Ave | 0.0665 | 0.0705 | | 0.0424 | 0.0400 | 6.0 | 20.0 |
| Chloroform | Ave | 0.3979 | 0.4367 | 0.2000 | 0.0219 | 0.0200 | 9.7 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3127 | 0.3465 | 0.1000 | 0.0222 | 0.0200 | 10.8 | 20.0 |
| Cyclohexane | Ave | 0.5900 | 0.6590 | 0.1000 | 0.0223 | 0.0200 | 11.7 | 20.0 |
| Carbon tetrachloride | Ave | 0.2541 | 0.2901 | 0.1000 | 0.0228 | 0.0200 | 14.2 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3331 | 0.3884 | | 0.0233 | 0.0200 | 16.6 | 20.0 |
| Benzene | Ave | 0.999 | 1.113 | 0.5000 | 0.0223 | 0.0200 | 11.4 | 20.0 |
| Isobutyl alcohol | Ave | 0.0071 | 0.0082 | | 0.573 | 0.500 | 14.6 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.2792 | 0.3100 | 0.1000 | 0.0222 | 0.0200 | 11.0 | 20.0 |
| n-Heptane | Ave | 0.0682 | 0.0773 | | 0.0227 | 0.0200 | 13.4 | 20.0 |
| Trichloroethene | Ave | 0.2336 | 0.2694 | 0.2000 | 0.0231 | 0.0200 | 15.3 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-196026/13  Calibration Date: 09/02/2015 18:07

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 144945.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.4826 | 0.5370 | 0.1000 | 0.0223 | 0.0200 | 11.3 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2493 | 0.2803 | 0.1000 | 0.0225 | 0.0200 | 12.5 | 20.0 |
| Dibromomethane | Ave | 0.1040 | 0.1205 | | 0.0232 | 0.0200 | 15.9 | 20.0 |
| 1,4-Dioxane | Ave | 0.0017 | 0.0019 | | 0.462 | 0.400 | 15.4 | 20.0 |
| Bromodichloromethane | Ave | 0.2508 | 0.2588 | 0.1000 | 0.0206 | 0.0200 | 3.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0933 | 0.1063 | | 0.0228 | 0.0200 | 14.0 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3090 | 0.3296 | 0.1500 | 0.0213 | 0.0200 | 6.7 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3102 | 0.3490 | 0.0500 | 0.0450 | 0.0400 | 12.5 | 20.0 |
| Toluene | Ave | 1.617 | 1.748 | 0.4000 | 0.0216 | 0.0200 | 8.1 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3680 | 0.4131 | 0.1000 | 0.0224 | 0.0200 | 12.2 | 20.0 |
| Ethyl methacrylate | Ave | 0.3012 | 0.3245 | | 0.0215 | 0.0200 | 7.7 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2333 | 0.2665 | 0.1000 | 0.0228 | 0.0200 | 14.2 | 20.0 |
| Tetrachloroethene | Ave | 0.2833 | 0.3198 | 0.2000 | 0.0226 | 0.0200 | 12.9 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4264 | 0.4668 | | 0.0219 | 0.0200 | 9.5 | 20.0 |
| 2-Hexanone | Ave | 0.2007 | 0.2198 | 0.0500 | 0.0438 | 0.0400 | 9.5 | 20.0 |
| Dibromochloromethane | Lin1 | | 0.2290 | | 0.0193 | 0.0200 | -3.6 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2138 | 0.2523 | | 0.0236 | 0.0200 | 18.0 | 20.0 |
| Chlorobenzene | Ave | 0.9298 | 1.009 | 0.3000 | 0.0217 | 0.0200 | 8.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.2902 | 0.3344 | | 0.0231 | 0.0200 | 15.3 | 20.0 |
| Ethylbenzene | Ave | 0.5296 | 0.5798 | | 0.0219 | 0.0200 | 9.5 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6468 | 0.6849 | | 0.0212 | 0.0200 | 5.9 | 20.0 |
| o-Xylene | Ave | 0.6346 | 0.7209 | | 0.0227 | 0.0200 | 13.6 | 20.0 |
| Styrene | Ave | 0.9556 | 0.9735 | 0.3000 | 0.0204 | 0.0200 | 1.9 | 20.0 |
| Bromoform | Lin1 | | 0.1061 | 0.1000 | 0.0173 | 0.0200 | -13.4 | 20.0 |
| Isopropylbenzene | Ave | 1.742 | 1.946 | 0.1000 | 0.0223 | 0.0200 | 11.7 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.5958 | 0.6826 | 0.3000 | 0.0229 | 0.0200 | 14.6 | 20.0 |
| Bromobenzene | Ave | 0.7042 | 0.7858 | | 0.0223 | 0.0200 | 11.6 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1763 | 0.2024 | | 0.0230 | 0.0200 | 14.8 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2152 | 0.2320 | | 0.0216 | 0.0200 | 7.8 | 20.0 |
| N-Propylbenzene | Ave | 0.9769 | 1.052 | | 0.0215 | 0.0200 | 7.7 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8014 | 0.8578 | | 0.0214 | 0.0200 | 7.0 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.102 | 3.417 | | 0.0220 | 0.0200 | 10.2 | 20.0 |
| 4-Chlorotoluene | Ave | 0.7776 | 0.8277 | | 0.0213 | 0.0200 | 6.5 | 20.0 |
| tert-Butylbenzene | Ave | 2.752 | 3.074 | | 0.0223 | 0.0200 | 11.7 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.081 | 3.372 | | 0.0219 | 0.0200 | 9.5 | 20.0 |
| sec-Butylbenzene | Ave | 4.194 | 4.751 | | 0.0227 | 0.0200 | 13.3 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.474 | 1.583 | 0.6000 | 0.0215 | 0.0200 | 7.4 | 20.0 |
| p-Isopropyltoluene | Ave | 3.242 | 3.585 | | 0.0221 | 0.0200 | 10.6 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.484 | 1.542 | 0.5000 | 0.0208 | 0.0200 | 3.9 | 20.0 |
| n-Butylbenzene | Ave | 3.148 | 3.388 | | 0.0215 | 0.0200 | 7.6 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.421 | 1.521 | 0.4000 | 0.0214 | 0.0200 | 7.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-196026/13 Calibration Date: 09/02/2015 18:07

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 144945.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0871 | 0.0879 | 0.0500 | 0.0202 | 0.0200 | 0.9 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.9733 | 0.9859 | 0.2000 | 0.0203 | 0.0200 | 1.3 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5396 | 0.6222 | | 0.0231 | 0.0200 | 15.3 | 20.0 |
| Naphthalene | Ave | 2.263 | 2.304 | | 0.0204 | 0.0200 | 1.8 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.9223 | 0.9287 | | 0.0201 | 0.0200 | 0.7 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2154 | 0.2054 | | 0.0477 | 0.0500 | -4.6 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2516 | 0.2368 | | 0.0471 | 0.0500 | -5.9 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.539 | 1.437 | | 0.0467 | 0.0500 | -6.6 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.147 | 1.045 | | 0.0455 | 0.0500 | -8.9 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206051/3 Calibration Date: 11/10/2015 10:47

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 146325.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2422 | 0.3028 | 0.1000 | 0.0250 | 0.0200 | 25.0* | 20.0 |
| Chloromethane | Ave | 0.3787 | 0.3469 | 0.1000 | 0.0183 | 0.0200 | -8.4 | 20.0 |
| Vinyl chloride | Ave | 0.2747 | 0.2618 | 0.1000 | 0.0191 | 0.0200 | -4.7 | 20.0 |
| Butadiene | Ave | 0.3003 | 0.2963 | | 0.0197 | 0.0200 | -1.3 | 20.0 |
| Bromomethane | Ave | 0.1383 | 0.1443 | 0.0500 | 0.0209 | 0.0200 | 4.3 | 20.0 |
| Chloroethane | Ave | 0.1448 | 0.1576 | 0.0500 | 0.0218 | 0.0200 | 8.9 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3245 | 0.3591 | | 0.0221 | 0.0200 | 10.7 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2852 | 0.3576 | 0.1000 | 0.0251 | 0.0200 | 25.4* | 20.0 |
| Ethyl ether | Ave | 0.1804 | 0.1677 | | 0.0465 | 0.0500 | -7.0 | 20.0 |
| Acrolein | Ave | 0.0180 | 0.0143 | | 0.199 | 0.250 | -20.5* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2217 | 0.2482 | 0.1000 | 0.0560 | 0.0500 | 12.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1625 | 0.1959 | 0.0500 | 0.0603 | 0.0500 | 20.6* | 20.0 |
| Acetone | Lin1 | | 0.0643 | 0.0500 | 0.0926 | 0.100 | -7.4 | 20.0 |
| Iodomethane | Ave | 0.3194 | 0.3367 | | 0.0527 | 0.0500 | 5.4 | 20.0 |
| Carbon disulfide | Ave | 0.6714 | 0.6672 | 0.1000 | 0.0497 | 0.0500 | -0.6 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1430 | 0.1423 | | 0.0497 | 0.0500 | -0.5 | 20.0 |
| Methyl acetate | Ave | 0.1567 | 0.1298 | 0.1000 | 0.207 | 0.250 | -17.2 | 20.0 |
| Methylene Chloride | Lin1 | | 0.2547 | 0.1000 | 0.0494 | 0.0500 | -1.2 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0198 | 0.0165 | | 0.415 | 0.500 | -17.1 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2515 | 0.2656 | 0.1000 | 0.0528 | 0.0500 | 5.6 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.5650 | 0.5586 | 0.1000 | 0.0494 | 0.0500 | -1.1 | 20.0 |
| Acrylonitrile | Ave | 0.0757 | 0.0653 | | 0.431 | 0.500 | -13.8 | 20.0 |
| Hexane | Ave | 0.0816 | 0.0754 | | 0.0462 | 0.0500 | -7.5 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4925 | 0.4861 | 0.1000 | 0.0493 | 0.0500 | -1.3 | 20.0 |
| Vinyl acetate | Lin1 | | 0.0338 | | 0.0562 | 0.0500 | 12.4 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2107 | 0.2815 | | 0.0668 | 0.0500 | 33.6* | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2654 | 0.2800 | 0.1000 | 0.0528 | 0.0500 | 5.5 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0962 | 0.0843 | 0.0500 | 0.0876 | 0.100 | -12.4 | 20.0 |
| Bromochloromethane | Ave | 0.0967 | 0.1095 | | 0.0566 | 0.0500 | 13.2 | 20.0 |
| Tetrahydrofuran | Ave | 0.0665 | 0.0555 | | 0.0835 | 0.100 | -16.5 | 20.0 |
| Chloroform | Ave | 0.3979 | 0.4268 | 0.2000 | 0.0536 | 0.0500 | 7.3 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3127 | 0.3856 | 0.1000 | 0.0616 | 0.0500 | 23.3* | 20.0 |
| Cyclohexane | Ave | 0.5900 | 0.5511 | 0.1000 | 0.0467 | 0.0500 | -6.6 | 20.0 |
| Carbon tetrachloride | Ave | 0.2541 | 0.3080 | 0.1000 | 0.0606 | 0.0500 | 21.2* | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3331 | 0.3407 | | 0.0511 | 0.0500 | 2.3 | 20.0 |
| Benzene | Ave | 0.999 | 1.002 | 0.5000 | 0.0502 | 0.0500 | 0.3 | 20.0 |
| Isobutyl alcohol | Ave | 0.0071 | 0.0058 | | 1.02 | 1.25 | -18.5 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.2792 | 0.2827 | 0.1000 | 0.0506 | 0.0500 | 1.2 | 20.0 |
| n-Heptane | Ave | 0.0682 | 0.0764 | | 0.0560 | 0.0500 | 12.1 | 20.0 |
| Trichloroethene | Ave | 0.2336 | 0.2579 | 0.2000 | 0.0552 | 0.0500 | 10.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206051/3 Calibration Date: 11/10/2015 10:47

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 146325.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.4826 | 0.5136 | 0.1000 | 0.0532 | 0.0500 | 6.4 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2493 | 0.2442 | 0.1000 | 0.0490 | 0.0500 | -2.0 | 20.0 |
| Dibromomethane | Ave | 0.1040 | 0.1112 | | 0.0535 | 0.0500 | 6.9 | 20.0 |
| 1,4-Dioxane | Ave | 0.0017 | 0.0017 | | 1.03 | 1.00 | 3.3 | 20.0 |
| Bromodichloromethane | Ave | 0.2508 | 0.2740 | 0.1000 | 0.0546 | 0.0500 | 9.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0933 | 0.0821 | | 0.0880 | 0.100 | -12.0 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3090 | 0.3112 | 0.1500 | 0.0504 | 0.0500 | 0.7 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3102 | 0.2676 | 0.0500 | 0.0863 | 0.100 | -13.7 | 20.0 |
| Toluene | Ave | 1.617 | 1.728 | 0.4000 | 0.0534 | 0.0500 | 6.9 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3680 | 0.3806 | 0.1000 | 0.0517 | 0.0500 | 3.4 | 20.0 |
| Ethyl methacrylate | Ave | 0.3012 | 0.2921 | | 0.0485 | 0.0500 | -3.0 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2333 | 0.2582 | 0.1000 | 0.0553 | 0.0500 | 10.6 | 20.0 |
| Tetrachloroethene | Ave | 0.2833 | 0.3043 | 0.2000 | 0.0537 | 0.0500 | 7.4 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4264 | 0.4318 | | 0.0506 | 0.0500 | 1.3 | 20.0 |
| 2-Hexanone | Ave | 0.2007 | 0.1769 | 0.0500 | 0.0881 | 0.100 | -11.9 | 20.0 |
| Dibromochloromethane | Lin1 | | 0.2579 | | 0.0520 | 0.0500 | 3.9 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2138 | 0.2359 | | 0.0552 | 0.0500 | 10.3 | 20.0 |
| Chlorobenzene | Ave | 0.9298 | 0.9905 | 0.3000 | 0.0533 | 0.0500 | 6.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.2902 | 0.3234 | | 0.0557 | 0.0500 | 11.5 | 20.0 |
| Ethylbenzene | Ave | 0.5296 | 0.5533 | | 0.0522 | 0.0500 | 4.5 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6468 | 0.6818 | | 0.0527 | 0.0500 | 5.4 | 20.0 |
| o-Xylene | Ave | 0.6346 | 0.6810 | | 0.0537 | 0.0500 | 7.3 | 20.0 |
| Styrene | Ave | 0.9556 | 1.032 | 0.3000 | 0.0540 | 0.0500 | 8.0 | 20.0 |
| Bromoform | Lin1 | | 0.1280 | 0.1000 | 0.0502 | 0.0500 | 0.3 | 20.0 |
| Isopropylbenzene | Ave | 1.742 | 1.907 | 0.1000 | 0.0547 | 0.0500 | 9.5 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.5958 | 0.6150 | 0.3000 | 0.0516 | 0.0500 | 3.2 | 20.0 |
| Bromobenzene | Ave | 0.7042 | 0.7856 | | 0.0558 | 0.0500 | 11.6 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1763 | 0.1912 | | 0.0542 | 0.0500 | 8.5 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2152 | 0.1899 | | 0.0441 | 0.0500 | -11.8 | 20.0 |
| N-Propylbenzene | Ave | 0.9769 | 1.061 | | 0.0543 | 0.0500 | 8.6 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8014 | 0.8759 | | 0.0546 | 0.0500 | 9.3 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.102 | 3.395 | | 0.0547 | 0.0500 | 9.5 | 20.0 |
| 4-Chlorotoluene | Ave | 0.7776 | 0.8372 | | 0.0538 | 0.0500 | 7.7 | 20.0 |
| tert-Butylbenzene | Ave | 2.752 | 3.033 | | 0.0551 | 0.0500 | 10.2 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.081 | 3.382 | | 0.0549 | 0.0500 | 9.8 | 20.0 |
| sec-Butylbenzene | Ave | 4.194 | 4.534 | | 0.0540 | 0.0500 | 8.1 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.474 | 1.485 | 0.6000 | 0.0504 | 0.0500 | 0.7 | 20.0 |
| p-Isopropyltoluene | Ave | 3.242 | 3.675 | | 0.0567 | 0.0500 | 13.4 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.484 | 1.475 | 0.5000 | 0.0497 | 0.0500 | -0.6 | 20.0 |
| n-Butylbenzene | Ave | 3.148 | 3.476 | | 0.0552 | 0.0500 | 10.4 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.421 | 1.407 | 0.4000 | 0.0495 | 0.0500 | -1.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206051/3  Calibration Date: 11/10/2015 10:47

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146325.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0871 | 0.0936 | 0.0500 | 0.0537 | 0.0500 | 7.4 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.9733 | 1.040 | 0.2000 | 0.0534 | 0.0500 | 6.8 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5396 | 0.6133 | | 0.0568 | 0.0500 | 13.7 | 20.0 |
| Naphthalene | Ave | 2.263 | 2.099 | | 0.0464 | 0.0500 | -7.2 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.9223 | 0.9394 | | 0.0509 | 0.0500 | 1.9 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2154 | 0.2257 | | 0.0524 | 0.0500 | 4.8 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2516 | 0.2552 | | 0.0507 | 0.0500 | 1.5 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.539 | 1.542 | | 0.0501 | 0.0500 | 0.2 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.147 | 1.068 | | 0.0465 | 0.0500 | -7.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206051/4  Calibration Date: 11/10/2015 11:08

Instrument ID: A3UX14  Calib Start Date: 08/13/2015 17:50

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 08/13/2015 20:42

Lab File ID: 146326.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0224 | 0.0216 | | 0.481 | 0.500 | -3.7 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.4341 | 0.4643 | | 0.0535 | 0.0500 | 7.0 | 20.0 |
| Diisopropyl ether | Ave | 0.2079 | 0.2164 | | 0.0520 | 0.0500 | 4.1 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8262 | 0.7658 | | 0.0463 | 0.0500 | -7.3 | 20.0 |
| Ethyl acetate | Ave | 0.1606 | 0.1374 | | 0.0855 | 0.100 | -14.5 | 20.0 |
| Propionitrile | Ave | 0.0278 | 0.0233 | | 0.420 | 0.500 | -16.0 | 20.0 |
| Methacrylonitrile | Ave | 0.1070 | 0.1039 | | 0.486 | 0.500 | -2.9 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.7020 | 0.6252 | | 0.0445 | 0.0500 | -10.9 | 20.0 |
| n-Butanol | Ave | 0.0105 | 0.0071 | | 0.846 | 1.25 | -32.4* | 20.0 |
| Methyl methacrylate | Ave | 0.1432 | 0.1329 | | 0.0928 | 0.100 | -7.2 | 20.0 |
| 2-Nitropropane | Ave | 0.0224 | 0.0235 | | 0.105 | 0.100 | 4.8 | 20.0 |
| 1-Chlorohexane | Ave | 0.4504 | 0.5066 | | 0.0562 | 0.0500 | 12.5 | 20.0 |
| Cyclohexanone | Ave | 0.0123 | 0.0082 | | 0.334 | 0.500 | -33.3* | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.034 | 3.484 | | 0.0574 | 0.0500 | 14.8 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.096 | 1.298 | | 0.0592 | 0.0500 | 18.4 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.220 | 1.397 | | 0.115 | 0.100 | 14.5 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206267/3 Calibration Date: 11/11/2015 10:39

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 146358.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2422 | 0.2811 | 0.1000 | 0.0232 | 0.0200 | 16.0 | 20.0 |
| Chloromethane | Ave | 0.3787 | 0.3232 | 0.1000 | 0.0171 | 0.0200 | -14.6 | 20.0 |
| Vinyl chloride | Ave | 0.2747 | 0.2526 | 0.1000 | 0.0184 | 0.0200 | -8.0 | 20.0 |
| Butadiene | Ave | 0.3003 | 0.2846 | | 0.0190 | 0.0200 | -5.2 | 20.0 |
| Bromomethane | Ave | 0.1383 | 0.1381 | 0.0500 | 0.0200 | 0.0200 | -0.2 | 20.0 |
| Chloroethane | Ave | 0.1448 | 0.1525 | 0.0500 | 0.0211 | 0.0200 | 5.3 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3245 | 0.3469 | | 0.0214 | 0.0200 | 6.9 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2852 | 0.3317 | 0.1000 | 0.0233 | 0.0200 | 16.3 | 20.0 |
| Ethyl ether | Ave | 0.1804 | 0.1573 | | 0.0436 | 0.0500 | -12.8 | 20.0 |
| Acrolein | Ave | 0.0180 | 0.0124 | | 0.172 | 0.250 | -31.2* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2217 | 0.2132 | 0.1000 | 0.0481 | 0.0500 | -3.8 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1625 | 0.1832 | 0.0500 | 0.0564 | 0.0500 | 12.8 | 20.0 |
| Acetone | Lin1 | | 0.0558 | 0.0500 | 0.0790 | 0.100 | -21.0* | 20.0 |
| Iodomethane | Ave | 0.3194 | 0.3184 | | 0.0499 | 0.0500 | -0.3 | 20.0 |
| Carbon disulfide | Ave | 0.6714 | 0.6540 | 0.1000 | 0.0487 | 0.0500 | -2.6 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1430 | 0.1337 | | 0.0467 | 0.0500 | -6.5 | 20.0 |
| Methyl acetate | Ave | 0.1567 | 0.1142 | 0.1000 | 0.182 | 0.250 | -27.1* | 20.0 |
| Methylene Chloride | Lin1 | | 0.2445 | 0.1000 | 0.0472 | 0.0500 | -5.7 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0198 | 0.0143 | | 0.359 | 0.500 | -28.2* | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2515 | 0.2557 | 0.1000 | 0.0508 | 0.0500 | 1.7 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.5650 | 0.5068 | 0.1000 | 0.0448 | 0.0500 | -10.3 | 20.0 |
| Acrylonitrile | Ave | 0.0757 | 0.0582 | | 0.384 | 0.500 | -23.1* | 20.0 |
| Hexane | Ave | 0.0816 | 0.0688 | | 0.0422 | 0.0500 | -15.6 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4925 | 0.4603 | 0.1000 | 0.0467 | 0.0500 | -6.5 | 20.0 |
| Vinyl acetate | Lin1 | | 0.0268 | | 0.0448 | 0.0500 | -10.4 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2107 | 0.2606 | | 0.0618 | 0.0500 | 23.7* | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2654 | 0.2665 | 0.1000 | 0.0502 | 0.0500 | 0.4 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0962 | 0.0686 | 0.0500 | 0.0713 | 0.100 | -28.7* | 20.0 |
| Bromochloromethane | Ave | 0.0967 | 0.1012 | | 0.0523 | 0.0500 | 4.6 | 20.0 |
| Tetrahydrofuran | Ave | 0.0665 | 0.0474 | | 0.0713 | 0.100 | -28.7* | 20.0 |
| Chloroform | Ave | 0.3979 | 0.4075 | 0.2000 | 0.0512 | 0.0500 | 2.4 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3127 | 0.3691 | 0.1000 | 0.0590 | 0.0500 | 18.0 | 20.0 |
| Cyclohexane | Ave | 0.5900 | 0.5303 | 0.1000 | 0.0449 | 0.0500 | -10.1 | 20.0 |
| Carbon tetrachloride | Ave | 0.2541 | 0.2972 | 0.1000 | 0.0585 | 0.0500 | 17.0 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3331 | 0.3194 | | 0.0479 | 0.0500 | -4.1 | 20.0 |
| Benzene | Ave | 0.999 | 0.9482 | 0.5000 | 0.0475 | 0.0500 | -5.1 | 20.0 |
| Isobutyl alcohol | Ave | 0.0071 | 0.0050 | | 0.873 | 1.25 | -30.2* | 20.0 |
| 1,2-Dichloroethane | Ave | 0.2792 | 0.2670 | 0.1000 | 0.0478 | 0.0500 | -4.4 | 20.0 |
| n-Heptane | Ave | 0.0682 | 0.0690 | | 0.0506 | 0.0500 | 1.2 | 20.0 |
| Trichloroethene | Ave | 0.2336 | 0.2428 | 0.2000 | 0.0520 | 0.0500 | 4.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206267/3  Calibration Date: 11/11/2015 10:39

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146358.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.4826 | 0.4880 | 0.1000 | 0.0506 | 0.0500 | 1.1 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2493 | 0.2230 | 0.1000 | 0.0447 | 0.0500 | -10.5 | 20.0 |
| Dibromomethane | Ave | 0.1040 | 0.1015 | | 0.0488 | 0.0500 | -2.4 | 20.0 |
| 1,4-Dioxane | Ave | 0.0017 | 0.0013 | | 0.789 | 1.00 | -21.1* | 20.0 |
| Bromodichloromethane | Ave | 0.2508 | 0.2514 | 0.1000 | 0.0501 | 0.0500 | 0.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0933 | 0.0673 | | 0.0721 | 0.100 | -27.9* | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3090 | 0.2735 | 0.1500 | 0.0443 | 0.0500 | -11.5 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3102 | 0.2327 | 0.0500 | 0.0750 | 0.100 | -25.0* | 20.0 |
| Toluene | Ave | 1.617 | 1.639 | 0.4000 | 0.0507 | 0.0500 | 1.4 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3680 | 0.3380 | 0.1000 | 0.0459 | 0.0500 | -8.2 | 20.0 |
| Ethyl methacrylate | Ave | 0.3012 | 0.2523 | | 0.0419 | 0.0500 | -16.2 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2333 | 0.2322 | 0.1000 | 0.0498 | 0.0500 | -0.5 | 20.0 |
| Tetrachloroethene | Ave | 0.2833 | 0.2858 | 0.2000 | 0.0505 | 0.0500 | 0.9 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4264 | 0.3916 | | 0.0459 | 0.0500 | -8.2 | 20.0 |
| 2-Hexanone | Ave | 0.2007 | 0.1471 | 0.0500 | 0.0733 | 0.100 | -26.7* | 20.0 |
| Dibromochloromethane | Lin1 | | 0.2357 | | 0.0476 | 0.0500 | -4.8 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2138 | 0.2082 | | 0.0487 | 0.0500 | -2.6 | 20.0 |
| Chlorobenzene | Ave | 0.9298 | 0.9062 | 0.3000 | 0.0487 | 0.0500 | -2.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.2902 | 0.3043 | | 0.0524 | 0.0500 | 4.9 | 20.0 |
| Ethylbenzene | Ave | 0.5296 | 0.5106 | | 0.0482 | 0.0500 | -3.6 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6468 | 0.6472 | | 0.0500 | 0.0500 | 0.0 | 20.0 |
| o-Xylene | Ave | 0.6346 | 0.6427 | | 0.0506 | 0.0500 | 1.3 | 20.0 |
| Styrene | Ave | 0.9556 | 0.9588 | 0.3000 | 0.0502 | 0.0500 | 0.3 | 20.0 |
| Bromoform | Lin1 | | 0.1101 | 0.1000 | 0.0433 | 0.0500 | -13.4 | 20.0 |
| Isopropylbenzene | Ave | 1.742 | 1.827 | 0.1000 | 0.0524 | 0.0500 | 4.9 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.5958 | 0.5674 | 0.3000 | 0.0476 | 0.0500 | -4.8 | 20.0 |
| Bromobenzene | Ave | 0.7042 | 0.7254 | | 0.0515 | 0.0500 | 3.0 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1763 | 0.1674 | | 0.0475 | 0.0500 | -5.0 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2152 | 0.1618 | | 0.0376 | 0.0500 | -24.8* | 20.0 |
| N-Propylbenzene | Ave | 0.9769 | 1.026 | | 0.0525 | 0.0500 | 5.0 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8014 | 0.8181 | | 0.0510 | 0.0500 | 2.1 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.102 | 3.275 | | 0.0528 | 0.0500 | 5.6 | 20.0 |
| 4-Chlorotoluene | Ave | 0.7776 | 0.7791 | | 0.0501 | 0.0500 | 0.2 | 20.0 |
| tert-Butylbenzene | Ave | 2.752 | 2.911 | | 0.0529 | 0.0500 | 5.8 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.081 | 3.208 | | 0.0521 | 0.0500 | 4.1 | 20.0 |
| sec-Butylbenzene | Ave | 4.194 | 4.345 | | 0.0518 | 0.0500 | 3.6 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.474 | 1.377 | 0.6000 | 0.0467 | 0.0500 | -6.6 | 20.0 |
| p-Isopropyltoluene | Ave | 3.242 | 3.491 | | 0.0538 | 0.0500 | 7.7 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.484 | 1.348 | 0.5000 | 0.0454 | 0.0500 | -9.2 | 20.0 |
| n-Butylbenzene | Ave | 3.148 | 3.277 | | 0.0520 | 0.0500 | 4.1 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.421 | 1.338 | 0.4000 | 0.0471 | 0.0500 | -5.8 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206267/3  Calibration Date: 11/11/2015 10:39

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146358.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0871 | 0.0827 | 0.0500 | 0.0475 | 0.0500 | -5.1 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.9733 | 0.9676 | 0.2000 | 0.0497 | 0.0500 | -0.6 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5396 | 0.5669 | | 0.0525 | 0.0500 | 5.1 | 20.0 |
| Naphthalene | Ave | 2.263 | 1.925 | | 0.0425 | 0.0500 | -14.9 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.9223 | 0.8905 | | 0.0483 | 0.0500 | -3.5 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2154 | 0.2264 | | 0.0526 | 0.0500 | 5.1 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2516 | 0.2478 | | 0.0493 | 0.0500 | -1.5 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.539 | 1.561 | | 0.0507 | 0.0500 | 1.5 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.147 | 1.030 | | 0.0449 | 0.0500 | -10.3 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206267/4  Calibration Date: 11/11/2015 11:00

Instrument ID: A3UX14  Calib Start Date: 08/13/2015 17:50

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 08/13/2015 20:42

Lab File ID: 146359.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0224 | 0.0212 | | 0.473 | 0.500 | -5.3 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.4341 | 0.4417 | | 0.0509 | 0.0500 | 1.7 | 20.0 |
| Diisopropyl ether | Ave | 0.2079 | 0.2137 | | 0.0514 | 0.0500 | 2.8 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8262 | 0.7386 | | 0.0447 | 0.0500 | -10.6 | 20.0 |
| Ethyl acetate | Ave | 0.1606 | 0.1311 | | 0.0817 | 0.100 | -18.3 | 20.0 |
| Propionitrile | Ave | 0.0278 | 0.0225 | | 0.405 | 0.500 | -18.9 | 20.0 |
| Methacrylonitrile | Ave | 0.1070 | 0.1032 | | 0.482 | 0.500 | -3.5 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.7020 | 0.6097 | | 0.0434 | 0.0500 | -13.1 | 20.0 |
| n-Butanol | Ave | 0.0105 | 0.0072 | | 0.856 | 1.25 | -31.5* | 20.0 |
| Methyl methacrylate | Ave | 0.1432 | 0.1282 | | 0.0895 | 0.100 | -10.5 | 20.0 |
| 2-Nitropropane | Ave | 0.0224 | 0.0219 | | 0.0976 | 0.100 | -2.4 | 20.0 |
| 1-Chlorohexane | Ave | 0.4504 | 0.4749 | | 0.0527 | 0.0500 | 5.5 | 20.0 |
| Cyclohexanone | Ave | 0.0123 | 0.0087 | | 0.353 | 0.500 | -29.3* | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.034 | 3.431 | | 0.0565 | 0.0500 | 13.1 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.096 | 1.271 | | 0.0580 | 0.0500 | 16.0 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.220 | 1.395 | | 0.114 | 0.100 | 14.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204067/14 Calibration Date: 10/28/2015 15:21

Instrument ID: A3UX15 Calib Start Date: 10/28/2015 13:05

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/28/2015 14:59

Lab File ID: UXC3830.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.4156 | 0.4225 | 0.1000 | 0.0102 | 0.0100 | 1.7 | 20.0 |
| Chloromethane | Ave | 0.3011 | 0.3110 | 0.1000 | 0.0103 | 0.0100 | 3.3 | 20.0 |
| Vinyl chloride | Ave | 0.3298 | 0.3307 | 0.1000 | 0.0100 | 0.0100 | 0.3 | 20.0 |
| Butadiene | Ave | 0.2931 | 0.2931 | | 0.0100 | 0.0100 | 0.0 | 20.0 |
| Bromomethane | Ave | 0.1790 | 0.1830 | 0.0500 | 0.0102 | 0.0100 | 2.2 | 20.0 |
| Chloroethane | Ave | 0.1902 | 0.1873 | 0.0500 | 0.00985 | 0.0100 | -1.5 | 20.0 |
| Dichlorofluoromethane | Ave | 0.5274 | 0.5397 | | 0.0102 | 0.0100 | 2.3 | 20.0 |
| Trichlorofluoromethane | Ave | 0.4637 | 0.4802 | 0.1000 | 0.0104 | 0.0100 | 3.5 | 20.0 |
| Ethyl ether | Ave | 0.2188 | 0.2222 | | 0.0102 | 0.0100 | 1.6 | 20.0 |
| Acrolein | Ave | 0.0206 | 0.0213 | | 0.0519 | 0.0500 | 3.7 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.3129 | 0.3193 | 0.1000 | 0.0102 | 0.0100 | 2.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2564 | 0.2516 | 0.0500 | 0.00981 | 0.0100 | -1.9 | 20.0 |
| Acetone | Lin1 | | 0.0720 | 0.0500 | 0.0179 | 0.0200 | -10.4 | 20.0 |
| Iodomethane | Ave | 0.4880 | 0.5249 | | 0.0108 | 0.0100 | 7.6 | 20.0 |
| Carbon disulfide | Ave | 0.8576 | 0.9130 | 0.1000 | 0.0106 | 0.0100 | 6.5 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1725 | 0.1700 | | 0.00986 | 0.0100 | -1.4 | 20.0 |
| Methyl acetate | Ave | 0.1509 | 0.1547 | 0.1000 | 0.0513 | 0.0500 | 2.5 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3809 | 0.1000 | 0.0108 | 0.0100 | 8.1 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0133 | 0.0127 | | 0.0952 | 0.100 | -4.8 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7657 | 0.8017 | 0.1000 | 0.0105 | 0.0100 | 4.7 | 20.0 |
| Acrylonitrile | Ave | 0.0817 | 0.0848 | | 0.104 | 0.100 | 3.7 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.3449 | 0.3632 | 0.1000 | 0.0105 | 0.0100 | 5.3 | 20.0 |
| Hexane | Ave | 0.0890 | 0.0954 | | 0.0107 | 0.0100 | 7.2 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.5976 | 0.6462 | 0.1000 | 0.0108 | 0.0100 | 8.1 | 20.0 |
| Vinyl acetate | Ave | 0.3243 | 0.3397 | | 0.0105 | 0.0100 | 4.7 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.3227 | 0.3284 | | 0.0102 | 0.0100 | 1.8 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3710 | 0.3792 | 0.1000 | 0.0102 | 0.0100 | 2.2 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0913 | 0.0904 | 0.0500 | 0.0198 | 0.0200 | -0.9 | 20.0 |
| Bromochloromethane | Ave | 0.1664 | 0.1732 | | 0.0104 | 0.0100 | 4.1 | 20.0 |
| Tetrahydrofuran | Ave | 0.0518 | 0.0521 | | 0.0201 | 0.0200 | 0.6 | 20.0 |
| Chloroform | Ave | 0.6432 | 0.6593 | 0.2000 | 0.0103 | 0.0100 | 2.5 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.4847 | 0.5200 | 0.1000 | 0.0107 | 0.0100 | 7.3 | 20.0 |
| Cyclohexane | Ave | 0.5639 | 0.5858 | 0.1000 | 0.0104 | 0.0100 | 3.9 | 20.0 |
| Carbon tetrachloride | Ave | 0.4107 | 0.4418 | 0.1000 | 0.0108 | 0.0100 | 7.6 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.4747 | 0.5316 | | 0.0112 | 0.0100 | 12.0 | 20.0 |
| Isobutyl alcohol | Ave | 0.0049 | 0.0048 | | 0.240 | 0.250 | -3.8 | 20.0 |
| Benzene | Ave | 1.413 | 1.418 | 0.5000 | 0.0100 | 0.0100 | 0.3 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.5438 | 0.5527 | 0.1000 | 0.0102 | 0.0100 | 1.6 | 20.0 |
| n-Heptane | Ave | 0.0834 | 0.0834 | | 0.0100 | 0.0100 | 0.0 | 20.0 |
| Trichloroethene | Ave | 0.3644 | 0.3846 | 0.2000 | 0.0106 | 0.0100 | 5.5 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204067/14 Calibration Date: 10/28/2015 15:21

Instrument ID: A3UX15 Calib Start Date: 10/28/2015 13:05

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/28/2015 14:59

Lab File ID: UXC3830.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.5417 | 0.5577 | 0.1000 | 0.0103 | 0.0100 | 2.9 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2981 | 0.3108 | 0.1000 | 0.0104 | 0.0100 | 4.3 | 20.0 |
| Dibromomethane | Ave | 0.1769 | 0.1801 | | 0.0102 | 0.0100 | 1.8 | 20.0 |
| 1,4-Dioxane | Ave | 0.0026 | 0.0022 | | 0.173 | 0.200 | -13.3 | 20.0 |
| Bromodichloromethane | Ave | 0.3932 | 0.3997 | 0.1000 | 0.0102 | 0.0100 | 1.7 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.1542 | 0.1535 | | 0.00995 | 0.0100 | -0.5 | 20.0 |
| cis-1,3-Dichloropropene | Lin1 | | 0.4542 | 0.1500 | 0.00976 | 0.0100 | -2.4 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.1647 | 0.1659 | 0.0500 | 0.0201 | 0.0200 | 0.7 | 20.0 |
| Toluene | Ave | 1.969 | 2.020 | 0.4000 | 0.0103 | 0.0100 | 2.6 | 20.0 |
| trans-1,3-Dichloropropene | Lin1 | | 0.4894 | 0.1000 | 0.00966 | 0.0100 | -3.4 | 20.0 |
| Ethyl methacrylate | Lin1 | | 0.3989 | | 0.00973 | 0.0100 | -2.7 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.3317 | 0.3345 | 0.1000 | 0.0101 | 0.0100 | 0.8 | 20.0 |
| Tetrachloroethene | Ave | 0.3574 | 0.3694 | 0.2000 | 0.0103 | 0.0100 | 3.4 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.6349 | 0.6136 | | 0.00966 | 0.0100 | -3.4 | 20.0 |
| 2-Hexanone | Ave | 0.1536 | 0.1464 | 0.0500 | 0.0191 | 0.0200 | -4.7 | 20.0 |
| Dibromochloromethane | Lin1 | | 0.3187 | | 0.00900 | 0.0100 | -10.0 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.3251 | 0.3340 | | 0.0103 | 0.0100 | 2.7 | 20.0 |
| Chlorobenzene | Ave | 1.266 | 1.274 | 0.3000 | 0.0101 | 0.0100 | 0.6 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3736 | 0.3910 | | 0.0105 | 0.0100 | 4.6 | 20.0 |
| Ethylbenzene | Ave | 0.6530 | 0.6716 | | 0.0103 | 0.0100 | 2.9 | 20.0 |
| m-Xylene & p-Xylene | Ave | 1.789 | 1.810 | | 0.0101 | 0.0100 | 1.2 | 20.0 |
| o-Xylene | Ave | 0.7816 | 0.8009 | | 0.0102 | 0.0100 | 2.5 | 20.0 |
| Styrene | Ave | 1.363 | 1.339 | 0.3000 | 0.00982 | 0.0100 | -1.8 | 20.0 |
| Bromoform | Lin1 | | 0.1536 | 0.1000 | 0.00884 | 0.0100 | -11.6 | 20.0 |
| Isopropylbenzene | Ave | 2.107 | 2.197 | 0.1000 | 0.0104 | 0.0100 | 4.3 | 20.0 |
| Bromobenzene | Ave | 0.9352 | 0.9714 | | 0.0104 | 0.0100 | 3.9 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.6716 | 0.6952 | 0.3000 | 0.0104 | 0.0100 | 3.5 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.2436 | 0.2556 | | 0.0105 | 0.0100 | 4.9 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.1999 | 0.2133 | | 0.0107 | 0.0100 | 6.7 | 20.0 |
| N-Propylbenzene | Ave | 1.009 | 1.078 | | 0.0107 | 0.0100 | 6.8 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8945 | 0.9279 | | 0.0104 | 0.0100 | 3.7 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.344 | 3.567 | | 0.0107 | 0.0100 | 6.7 | 20.0 |
| 4-Chlorotoluene | Ave | 3.463 | 3.593 | | 0.0104 | 0.0100 | 3.7 | 20.0 |
| tert-Butylbenzene | Ave | 2.784 | 2.904 | | 0.0104 | 0.0100 | 4.3 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.455 | 3.643 | | 0.0105 | 0.0100 | 5.4 | 20.0 |
| sec-Butylbenzene | Ave | 3.779 | 3.924 | | 0.0104 | 0.0100 | 3.8 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.848 | 1.859 | 0.6000 | 0.0101 | 0.0100 | 0.6 | 20.0 |
| p-Isopropyltoluene | Ave | 3.438 | 3.575 | | 0.0104 | 0.0100 | 4.0 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.910 | 1.909 | 0.5000 | 0.00999 | 0.0100 | -0.0 | 20.0 |
| n-Butylbenzene | Ave | 2.836 | 2.887 | | 0.0102 | 0.0100 | 1.8 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.764 | 1.756 | 0.4000 | 0.00996 | 0.0100 | -0.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204067/14  Calibration Date: 10/28/2015 15:21

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 13:05

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 14:59

Lab File ID: UXC3830.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Lin1 | | 0.0997 | 0.0500 | 0.00916 | 0.0100 | -8.4 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 1.161 | 1.106 | 0.2000 | 0.00952 | 0.0100 | -4.8 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5590 | 0.5093 | | 0.00911 | 0.0100 | -8.9 | 20.0 |
| Naphthalene | Ave | 2.480 | 2.519 | | 0.0102 | 0.0100 | 1.6 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 1.053 | 1.024 | | 0.00972 | 0.0100 | -2.8 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.3239 | 0.3161 | | 0.00715 | 0.00733 | -2.4 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.4605 | 0.4381 | | 0.00697 | 0.00733 | -4.9 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.768 | 1.697 | | 0.00704 | 0.00733 | -4.0 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 0.6700 | 0.6795 | | 0.00744 | 0.00733 | 1.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204067/15  Calibration Date: 10/28/2015 18:00

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 15:44

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 17:37

Lab File ID: UXC3837.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0228 | 0.0226 | | 0.0993 | 0.100 | -0.7 | 20.0 |
| Diisopropyl ether | Ave | 0.2803 | 0.2896 | | 0.0103 | 0.0100 | 3.3 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.5186 | 0.5418 | | 0.0104 | 0.0100 | 4.5 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8119 | 0.8342 | | 0.0103 | 0.0100 | 2.7 | 20.0 |
| Ethyl acetate | Ave | 0.1644 | 0.1655 | | 0.0201 | 0.0200 | 0.7 | 20.0 |
| Propionitrile | Ave | 0.0293 | 0.0301 | | 0.103 | 0.100 | 2.7 | 20.0 |
| Methacrylonitrile | Ave | 0.1324 | 0.1374 | | 0.104 | 0.100 | 3.8 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.6462 | 0.6564 | | 0.0102 | 0.0100 | 1.6 | 20.0 |
| n-Butanol | Ave | 0.0035 | 0.0031 | | 0.219 | 0.250 | -12.3 | 20.0 |
| Methyl methacrylate | Ave | 0.1803 | 0.1875 | | 0.0208 | 0.0200 | 4.0 | 20.0 |
| 2-Nitropropane | Ave | 0.0274 | 0.0237 | | 0.0173 | 0.0200 | -13.7 | 20.0 |
| 1-Chlorohexane | Ave | 0.5138 | 0.5158 | | 0.0100 | 0.0100 | 0.4 | 20.0 |
| Cyclohexanone | Ave | 0.0128 | 0.0120 | | 0.0933 | 0.100 | -6.7 | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.502 | 3.574 | | 0.0102 | 0.0100 | 2.0 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.274 | 1.242 | | 0.00974 | 0.0100 | -2.6 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.424 | 1.167 | | 0.0164 | 0.0200 | -18.1 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206461/2 Calibration Date: 11/12/2015 10:41

Instrument ID: A3UX15 Calib Start Date: 10/28/2015 13:05

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/28/2015 14:59

Lab File ID: UXC4392.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.4156 | 0.4748 | 0.1000 | 0.0114 | 0.0100 | 14.3 | 20.0 |
| Chloromethane | Ave | 0.3011 | 0.2531 | 0.1000 | 0.00841 | 0.0100 | -15.9 | 20.0 |
| Vinyl chloride | Ave | 0.3298 | 0.2842 | 0.1000 | 0.00862 | 0.0100 | -13.8 | 20.0 |
| Butadiene | Ave | 0.2931 | 0.2523 | | 0.00861 | 0.0100 | -13.9 | 20.0 |
| Bromomethane | Ave | 0.1790 | 0.1968 | 0.0500 | 0.0110 | 0.0100 | 10.0 | 20.0 |
| Chloroethane | Ave | 0.1902 | 0.1836 | 0.0500 | 0.00965 | 0.0100 | -3.5 | 20.0 |
| Dichlorofluoromethane | Ave | 0.5274 | 0.5573 | | 0.0106 | 0.0100 | 5.7 | 20.0 |
| Trichlorofluoromethane | Ave | 0.4637 | 0.4470 | 0.1000 | 0.00964 | 0.0100 | -3.6 | 20.0 |
| Ethyl ether | Ave | 0.2188 | 0.1979 | | 0.00905 | 0.0100 | -9.5 | 20.0 |
| Acrolein | Ave | 0.0206 | 0.0157 | | 0.0381 | 0.0500 | -23.9* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.3129 | 0.3099 | 0.1000 | 0.00990 | 0.0100 | -1.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2564 | 0.2706 | 0.0500 | 0.0106 | 0.0100 | 5.5 | 20.0 |
| Acetone | Lin1 | | 0.0764 | 0.0500 | 0.0191 | 0.0200 | -4.3 | 20.0 |
| Iodomethane | Ave | 0.4880 | 0.5373 | | 0.0110 | 0.0100 | 10.1 | 20.0 |
| Carbon disulfide | Ave | 0.8576 | 0.8155 | 0.1000 | 0.00951 | 0.0100 | -4.9 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1725 | 0.1609 | | 0.00933 | 0.0100 | -6.7 | 20.0 |
| Methyl acetate | Ave | 0.1509 | 0.1291 | 0.1000 | 0.0428 | 0.0500 | -14.4 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3455 | 0.1000 | 0.00970 | 0.0100 | -3.0 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0133 | 0.0120 | | 0.0900 | 0.100 | -10.0 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7657 | 0.7865 | 0.1000 | 0.0103 | 0.0100 | 2.7 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.3449 | 0.3440 | 0.1000 | 0.00998 | 0.0100 | -0.2 | 20.0 |
| Acrylonitrile | Ave | 0.0817 | 0.0722 | | 0.0883 | 0.100 | -11.7 | 20.0 |
| Hexane | Ave | 0.0890 | 0.0821 | | 0.00922 | 0.0100 | -7.8 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.5976 | 0.5650 | 0.1000 | 0.00945 | 0.0100 | -5.5 | 20.0 |
| Vinyl acetate | Ave | 0.3243 | 0.3339 | | 0.0103 | 0.0100 | 2.9 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.3227 | 0.3420 | | 0.0106 | 0.0100 | 6.0 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3710 | 0.3555 | 0.1000 | 0.00958 | 0.0100 | -4.2 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0913 | 0.0782 | 0.0500 | 0.0171 | 0.0200 | -14.3 | 20.0 |
| Bromochloromethane | Ave | 0.1664 | 0.1838 | | 0.0110 | 0.0100 | 10.5 | 20.0 |
| Tetrahydrofuran | Ave | 0.0518 | 0.0467 | | 0.0181 | 0.0200 | -9.7 | 20.0 |
| Chloroform | Ave | 0.6432 | 0.6727 | 0.2000 | 0.0105 | 0.0100 | 4.6 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.4847 | 0.5235 | 0.1000 | 0.0108 | 0.0100 | 8.0 | 20.0 |
| Cyclohexane | Ave | 0.5639 | 0.4901 | 0.1000 | 0.00869 | 0.0100 | -13.1 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.4747 | 0.4704 | | 0.00991 | 0.0100 | -0.9 | 20.0 |
| Carbon tetrachloride | Ave | 0.4107 | 0.4780 | 0.1000 | 0.0116 | 0.0100 | 16.4 | 20.0 |
| Isobutyl alcohol | Ave | 0.0049 | 0.0045 | | 0.228 | 0.250 | -8.9 | 20.0 |
| Benzene | Ave | 1.413 | 1.334 | 0.5000 | 0.00944 | 0.0100 | -5.6 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.5438 | 0.5731 | 0.1000 | 0.0105 | 0.0100 | 5.4 | 20.0 |
| n-Heptane | Ave | 0.0834 | 0.0828 | | 0.00993 | 0.0100 | -0.7 | 20.0 |
| Trichloroethene | Ave | 0.3644 | 0.3950 | 0.2000 | 0.0108 | 0.0100 | 8.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206461/2  Calibration Date: 11/12/2015 10:41

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 13:05

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 14:59

Lab File ID: UXC4392.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.5417 | 0.4958 | 0.1000 | 0.00915 | 0.0100 | -8.5 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2981 | 0.2720 | 0.1000 | 0.00913 | 0.0100 | -8.7 | 20.0 |
| Dibromomethane | Ave | 0.1769 | 0.1852 | | 0.0105 | 0.0100 | 4.7 | 20.0 |
| 1,4-Dioxane | Ave | 0.0026 | 0.0024 | | 0.187 | 0.200 | -6.5 | 20.0 |
| Bromodichloromethane | Ave | 0.3932 | 0.4333 | 0.1000 | 0.0110 | 0.0100 | 10.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.1542 | 0.1349 | | 0.0175 | 0.0200 | -12.5 | 20.0 |
| cis-1,3-Dichloropropene | Lin1 | | 0.4161 | 0.1500 | 0.00898 | 0.0100 | -10.2 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.1647 | 0.1558 | 0.0500 | 0.0189 | 0.0200 | -5.4 | 20.0 |
| Toluene | Ave | 1.969 | 1.768 | 0.4000 | 0.00898 | 0.0100 | -10.2 | 20.0 |
| trans-1,3-Dichloropropene | Lin1 | | 0.4223 | 0.1000 | 0.00842 | 0.0100 | -15.8 | 20.0 |
| Ethyl methacrylate | Lin1 | | 0.3389 | | 0.00834 | 0.0100 | -16.6 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.3317 | 0.3105 | 0.1000 | 0.00936 | 0.0100 | -6.4 | 20.0 |
| Tetrachloroethene | Ave | 0.3574 | 0.3563 | 0.2000 | 0.00997 | 0.0100 | -0.3 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.6349 | 0.5496 | | 0.00866 | 0.0100 | -13.4 | 20.0 |
| 2-Hexanone | Ave | 0.1536 | 0.1235 | 0.0500 | 0.0161 | 0.0200 | -19.6 | 20.0 |
| Dibromochloromethane | Lin1 | | 0.3343 | | 0.00941 | 0.0100 | -5.9 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.3251 | 0.3105 | | 0.00955 | 0.0100 | -4.5 | 20.0 |
| Chlorobenzene | Ave | 1.266 | 1.196 | 0.3000 | 0.00944 | 0.0100 | -5.6 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3736 | 0.4061 | | 0.0109 | 0.0100 | 8.7 | 20.0 |
| Ethylbenzene | Ave | 0.6530 | 0.6239 | | 0.00956 | 0.0100 | -4.4 | 20.0 |
| m-Xylene & p-Xylene | Ave | 1.789 | 1.688 | | 0.00943 | 0.0100 | -5.7 | 20.0 |
| o-Xylene | Ave | 0.7816 | 0.7445 | | 0.00952 | 0.0100 | -4.8 | 20.0 |
| Styrene | Ave | 1.363 | 1.327 | 0.3000 | 0.00974 | 0.0100 | -2.6 | 20.0 |
| Bromoform | Lin1 | | 0.1723 | 0.1000 | 0.00983 | 0.0100 | -1.7 | 20.0 |
| Isopropylbenzene | Ave | 2.107 | 2.063 | 0.1000 | 0.00979 | 0.0100 | -2.1 | 20.0 |
| Bromobenzene | Ave | 0.9352 | 0.8652 | | 0.00925 | 0.0100 | -7.5 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.6716 | 0.5654 | 0.3000 | 0.00842 | 0.0100 | -15.8 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.2436 | 0.2253 | | 0.00925 | 0.0100 | -7.5 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.1999 | 0.1452 | | 0.00726 | 0.0100 | -27.4* | 20.0 |
| N-Propylbenzene | Ave | 1.009 | 0.8864 | | 0.00878 | 0.0100 | -12.2 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8945 | 0.7988 | | 0.00893 | 0.0100 | -10.7 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.344 | 2.940 | | 0.00879 | 0.0100 | -12.1 | 20.0 |
| 4-Chlorotoluene | Ave | 3.463 | 2.986 | | 0.00862 | 0.0100 | -13.8 | 20.0 |
| tert-Butylbenzene | Ave | 2.784 | 2.450 | | 0.00880 | 0.0100 | -12.0 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.455 | 3.098 | | 0.00897 | 0.0100 | -10.3 | 20.0 |
| sec-Butylbenzene | Ave | 3.779 | 3.234 | | 0.00856 | 0.0100 | -14.4 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.848 | 1.686 | 0.6000 | 0.00912 | 0.0100 | -8.8 | 20.0 |
| p-Isopropyltoluene | Ave | 3.438 | 3.118 | | 0.00907 | 0.0100 | -9.3 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.910 | 1.691 | 0.5000 | 0.00885 | 0.0100 | -11.5 | 20.0 |
| n-Butylbenzene | Ave | 2.836 | 2.292 | | 0.00808 | 0.0100 | -19.2 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.764 | 1.567 | 0.4000 | 0.00889 | 0.0100 | -11.1 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206461/2  Calibration Date: 11/12/2015 10:41

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 13:05

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 14:59

Lab File ID: UXC4392.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Lin1 | | 0.0997 | 0.0500 | 0.00917 | 0.0100 | -8.3 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 1.161 | 1.004 | 0.2000 | 0.00864 | 0.0100 | -13.6 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5590 | 0.4334 | | 0.00775 | 0.0100 | -22.5* | 20.0 |
| Naphthalene | Ave | 2.480 | 2.098 | | 0.00846 | 0.0100 | -15.4 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 1.053 | 0.8905 | | 0.00846 | 0.0100 | -15.4 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.3239 | 0.3247 | | 0.00735 | 0.00733 | 0.2 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.4605 | 0.4500 | | 0.00716 | 0.00733 | -2.3 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.768 | 1.546 | | 0.00641 | 0.00733 | -12.5 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 0.6700 | 0.6842 | | 0.00749 | 0.00733 | 2.1 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206461/3  Calibration Date: 11/12/2015 11:26

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 15:44

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 17:37

Lab File ID: UXC4394.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0228 | 0.0075 | | 0.0330 | 0.100 | -67.0* | 20.0 |
| Diisopropyl ether | Ave | 0.2803 | 0.2793 | | 0.00996 | 0.0100 | -0.4 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.5186 | 0.4985 | | 0.00961 | 0.0100 | -3.9 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8119 | 0.7747 | | 0.00954 | 0.0100 | -4.6 | 20.0 |
| Ethyl acetate | Ave | 0.1644 | 0.1216 | | 0.0148 | 0.0200 | -26.0* | 20.0 |
| Propionitrile | Ave | 0.0293 | 0.0194 | | 0.0663 | 0.100 | -33.7* | 20.0 |
| Methacrylonitrile | Ave | 0.1324 | 0.1064 | | 0.0804 | 0.100 | -19.6 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.6462 | 0.6591 | | 0.0102 | 0.0100 | 2.0 | 20.0 |
| n-Butanol | Ave | 0.0035 | 0.0016 | | 0.115 | 0.250 | -54.1* | 20.0 |
| Methyl methacrylate | Ave | 0.1803 | 0.1522 | | 0.0169 | 0.0200 | -15.6 | 20.0 |
| 2-Nitropropane | Ave | 0.0274 | 0.0252 | | 0.0184 | 0.0200 | -8.2 | 20.0 |
| 1-Chlorohexane | Ave | 0.5138 | 0.4605 | | 0.00896 | 0.0100 | -10.4 | 20.0 |
| Cyclohexanone | Ave | 0.0128 | 0.0048 | | 0.0378 | 0.100 | -62.2* | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.502 | 3.168 | | 0.00905 | 0.0100 | -9.5 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.274 | 1.080 | | 0.00847 | 0.0100 | -15.3 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.424 | 0.9190 | | 0.0129 | 0.0200 | -35.5* | 20.0 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206089/1-A

Matrix: Solid | Lab File ID: 146331.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5.0(g) | Date Analyzed: 11/10/2015 12:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.80 | U | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.78 | U | 5.0 | 0.78 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 2.83 | J | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.90 | U | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206089/1-A

Matrix: Solid | Lab File ID: 146331.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5.0(g) | Date Analyzed: 11/10/2015 12:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.79 | U | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.71 | U | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.7 | U | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.74 | U | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.35 | U | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.91 | U | 5.0 | 0.91 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.80 | U | 5.0 | 0.80 |
| 108-88-3 | Toluene | 0.27 | U | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206089/1-A

Matrix: Solid | Lab File ID: 146331.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5.0(g) | Date Analyzed: 11/10/2015 12:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 91 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 88 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 82 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 91 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206267/6

Matrix: Solid | Lab File ID: 146361.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/11/2015 11:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.80 | U | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.78 | U | 5.0 | 0.78 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 3.14 | J | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.90 | U | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206267/6

Matrix: Solid | Lab File ID: 146361.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/11/2015 11:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.79 | U | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.71 | U | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.7 | U | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.74 | U | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.350 | J | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.91 | U | 5.0 | 0.91 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.80 | U | 5.0 | 0.80 |
| 108-88-3 | Toluene | 0.369 | J | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206267/6

Matrix: Solid | Lab File ID: 146361.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/11/2015 11:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 101 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 90 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206461/6

Matrix: Water | Lab File ID: UXC4396.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:11

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.28 | U | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.44 | U | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.22 | U | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.30 | U | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 0.45 | U | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.37 | U | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.44 | U | 1.0 | 0.44 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.47 | U | 5.0 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.32 | U | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.41 | U | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.82 | U | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.25 | U | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 0.23 | U | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 0.25 | U | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.19 | U | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 0.19 | U | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.27 | U | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 0.26 | U | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 0.583 | J | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.59 | U | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 0.40 | U | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 0.48 | U | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 10 | 0.99 |
| 67-64-1 | Acetone | 0.94 | U | 10 | 0.94 |
| 71-43-2 | Benzene | 0.35 | U | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 0.35 | U | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 0.56 | U | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 0.44 | U | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.38 | U | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 0.43 | U | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 0.25 | U | 1.0 | 0.25 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206461/6

Matrix: Water | Lab File ID: UXC4396.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:11

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 1.0 | 0.43 |
| 75-00-3 | Chloroethane | 0.32 | U | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 0.25 | U | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 0.44 | U | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.26 | U | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.46 | U | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 0.42 | U | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.50 | U | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 4.9 | U | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 0.25 | U | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 0.35 | U | 1.0 | 0.35 |
| 108-20-3 | Diisopropyl ether | 0.50 | U | 10 | 0.50 |
| 98-82-8 | Isopropylbenzene | 0.35 | U | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 0.20 | U | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 0.33 | U | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.24 | U | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 0.45 | U | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 0.31 | U | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 0.40 | U | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 0.25 | U | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 0.48 | U | 1.0 | 0.48 |
| 100-42-5 | Styrene | 0.45 | U | 1.0 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.30 | U | 5.0 | 0.30 |
| 637-92-3 | Ethyl t-butyl ether | 0.23 | U | 5.0 | 0.23 |
| 127-18-4 | Tetrachloroethene | 0.31 | U | 1.0 | 0.31 |
| 108-88-3 | Toluene | 0.23 | U | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.30 | U | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.56 | U | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 0.22 | U | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 0.49 | U | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 0.41 | U | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 0.29 | U | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 0.52 | U | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 0.41 | U | 1.0 | 0.41 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-206461/6

Matrix: Water  Lab File ID: UXC4396.D

Analysis Method: 8260C  Date Collected:

Sample wt/vol: 5(mL)  Date Analyzed: 11/12/2015 12:11

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: DB-624  ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206461  Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.32 | U | 1.0 | 0.32 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 88 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 99 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 97 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 95 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206051/5

Matrix: Solid | Lab File ID: 146327.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/10/2015 11:30

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 20.4 | | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 22.0 | | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 19.6 | | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.9 | | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 18.5 | | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 19.1 | | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 20.3 | | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 19.5 | | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 20.3 | | 5.0 | 0.78 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 20.6 | | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 20.3 | | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 19.1 | | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 19.0 | | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 19.3 | | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 17.7 | | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 19.3 | | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 18.6 | | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 18.8 | | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 21.3 | | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 31.5 | | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 14.8 | J | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 20.2 | | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 31.3 | | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 20.0 | | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 20.7 | | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 30.7 | | 20 | 1.2 |
| 67-64-1 | Acetone | 32.8 | | 20 | 5.7 |
| 71-43-2 | Benzene | 18.5 | | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 21.3 | | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 16.3 | | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 18.8 | | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 15.1 | | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 21.9 | | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 19.9 | | 5.0 | 0.53 |
| 124-48-1 | Dibromochloromethane | 18.2 | | 5.0 | 0.37 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206051/5

Matrix: Solid | Lab File ID: 146327.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/10/2015 11:30

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 18.7 | | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 19.9 | | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 14.5 | | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 19.6 | | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 17.3 | | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 19.1 | | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 21.1 | | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 16.4 | | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 18.1 | | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 138 | J | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 19.8 | | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 21.4 | | 5.0 | 0.40 |
| 98-82-8 | Isopropylbenzene | 20.6 | | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 18.7 | | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 16.7 | | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 19.8 | | 10 | 0.42 |
| 91-20-3 | Naphthalene | 18.4 | | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 20.2 | | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 19.8 | | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 20.7 | | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 19.9 | | 5.0 | 0.91 |
| 100-42-5 | Styrene | 18.7 | | 5.0 | 0.40 |
| 127-18-4 | Tetrachloroethene | 19.6 | | 5.0 | 0.80 |
| 108-88-3 | Toluene | 20.5 | | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 20.1 | | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 19.3 | | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 21.1 | | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 22.3 | | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 16.9 | | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 15.8 | | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 40.5 | | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 20.1 | | 5.0 | 0.67 |
| 106-93-4 | 1,2-Dibromoethane | 19.7 | | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-206051/5

Matrix: Solid  Lab File ID: 146327.D

Analysis Method: 8260C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/10/2015 11:30

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: DB-624 ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206051  Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 94 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 95 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 92 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206267/5

Matrix: Solid | Lab File ID: 146360.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/11/2015 11:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 21.7 | | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 22.4 | | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 20.2 | | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 21.2 | | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 19.0 | | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 18.6 | | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 20.9 | | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 20.6 | | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 21.8 | | 5.0 | 0.78 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 21.3 | | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 21.1 | | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 20.9 | | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 20.6 | | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 20.6 | | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 19.1 | | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 20.2 | | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 20.5 | | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 20.4 | | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 21.9 | | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 33.7 | | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 15.3 | J | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 21.5 | | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 32.9 | | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 21.6 | | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 22.3 | | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 32.6 | | 20 | 1.2 |
| 67-64-1 | Acetone | 33.3 | | 20 | 5.7 |
| 71-43-2 | Benzene | 19.4 | | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 22.4 | | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 17.7 | | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 19.6 | | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 14.8 | | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 22.8 | | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 21.0 | | 5.0 | 0.53 |
| 124-48-1 | Dibromochloromethane | 19.4 | | 5.0 | 0.37 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206267/5

Matrix: Solid | Lab File ID: 146360.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/11/2015 11:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 19.9 | | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 21.0 | | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 15.6 | | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 20.1 | | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 18.4 | | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 21.4 | | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 22.2 | | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 20.4 | | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 19.2 | | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 143 | J | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 21.3 | | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 23.6 | | 5.0 | 0.40 |
| 98-82-8 | Isopropylbenzene | 21.2 | | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 18.5 | | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 16.9 | | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 21.3 | | 10 | 0.42 |
| 91-20-3 | Naphthalene | 18.4 | | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 21.8 | | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 21.7 | | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 21.7 | | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 20.9 | | 5.0 | 0.91 |
| 100-42-5 | Styrene | 20.0 | | 5.0 | 0.40 |
| 127-18-4 | Tetrachloroethene | 21.3 | | 5.0 | 0.80 |
| 108-88-3 | Toluene | 21.3 | | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 20.3 | | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 20.6 | | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 21.5 | | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 22.8 | | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 17.7 | | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 16.8 | | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 43.0 | | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 21.4 | | 5.0 | 0.67 |
| 106-93-4 | 1,2-Dibromoethane | 21.7 | | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-206267/5

Matrix: Solid  Lab File ID: 146360.D

Analysis Method: 8260C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/11/2015 11:22

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: DB-624 ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206267  Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 95 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 99 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 89 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206461/4

Matrix: Water | Lab File ID: UXC4393.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 11:03

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 11.8 | | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 11.1 | | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 9.38 | | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 9.57 | | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 9.43 | | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 8.73 | | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 10.4 | | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 8.64 | | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 11.3 | | 1.0 | 0.44 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 8.66 | | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 9.93 | | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 9.17 | | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 9.76 | | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 10.5 | | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 9.07 | | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 9.90 | | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 9.29 | | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 9.83 | | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 10.4 | | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 16.3 | | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 8.93 | J | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 10.4 | | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 17.6 | | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 9.89 | | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 9.42 | | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 18.1 | | 10 | 0.99 |
| 67-64-1 | Acetone | 13.6 | | 10 | 0.94 |
| 71-43-2 | Benzene | 9.23 | | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 10.7 | | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 10.7 | | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 11.0 | | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 7.86 | | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 11.6 | | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 10.2 | | 1.0 | 0.25 |
| 124-48-1 | Dibromochloromethane | 10.6 | | 1.0 | 0.43 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206461/4

Matrix: Water | Lab File ID: UXC4393.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 11:03

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 9.73 | | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 10.6 | | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 8.46 | | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 9.74 | | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 9.52 | | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 10.7 | | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 10.8 | | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 10.5 | | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 11.0 | | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 72.2 | | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 10.5 | | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 8.37 | | 1.0 | 0.35 |
| 98-82-8 | Isopropylbenzene | 10.4 | | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 10.1 | | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 9.34 | | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 10.0 | | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 8.44 | | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 8.71 | | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 10.2 | | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 10.4 | | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 9.27 | | 1.0 | 0.48 |
| 100-42-5 | Styrene | 9.85 | | 1.0 | 0.45 |
| 127-18-4 | Tetrachloroethene | 10.9 | | 1.0 | 0.31 |
| 108-88-3 | Toluene | 9.77 | | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 9.53 | | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 9.90 | | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 11.1 | | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 9.70 | | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 9.77 | | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 9.07 | | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 20.4 | | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 9.67 | | 1.0 | 0.41 |
| 106-93-4 | 1,2-Dibromoethane | 10.7 | | 1.0 | 0.32 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-206461/4

Matrix: Water  Lab File ID: UXC4393.D

Analysis Method: 8260C  Date Collected:

Sample wt/vol: 5(mL)  Date Analyzed: 11/12/2015 11:03

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: DB-624 ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206461  Units: ug/L

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 96 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 103 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 103 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 99 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 MS | Lab Sample ID: 240-57597-5 MS

Matrix: Solid | Lab File ID: 146341.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.17(g) | Date Analyzed: 11/10/2015 16:30

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 23.0 | | 5.7 | 0.57 |
| 71-55-6 | 1,1,1-Trichloroethane | 24.1 | | 5.7 | 0.74 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 26.6 | | 5.7 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 23.5 | | 5.7 | 0.48 |
| 75-34-3 | 1,1-Dichloroethane | 21.1 | | 5.7 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 19.8 | | 5.7 | 0.91 |
| 563-58-6 | 1,1-Dichloropropene | 22.7 | | 5.7 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 14.1 | | 5.7 | 0.57 |
| 96-18-4 | 1,2,3-Trichloropropane | 23.8 | | 5.7 | 0.89 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 1.81 | J | 5.7 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 14.2 | | 5.7 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 18.7 | | 5.7 | 0.66 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 22.2 | | 11 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 18.1 | | 5.7 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 22.0 | | 5.7 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 20.8 | | 5.7 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 18.3 | | 5.7 | 0.79 |
| 142-28-9 | 1,3-Dichloropropane | 22.5 | | 5.7 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 17.8 | | 5.7 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 24.0 | | 5.7 | 0.50 |
| 78-93-3 | 2-Butanone (MEK) | 36.6 | | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 16.4 | J | 57 | 0.73 |
| 95-49-8 | 2-Chlorotoluene | 20.8 | | 5.7 | 0.34 |
| 591-78-6 | 2-Hexanone | 52.2 | | 23 | 0.78 |
| 106-43-4 | 4-Chlorotoluene | 19.4 | | 5.7 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 19.0 | | 5.7 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 150 | | 23 | 1.4 |
| 67-64-1 | Acetone | 47.7 | | 23 | 6.5 |
| 71-43-2 | Benzene | 21.1 | | 5.7 | 1.0 |
| 108-86-1 | Bromobenzene | 21.7 | | 5.7 | 0.38 |
| 75-25-2 | Bromoform | 18.7 | | 5.7 | 0.27 |
| 74-83-9 | Bromomethane | 19.8 | | 5.7 | 0.48 |
| 75-15-0 | Carbon disulfide | 16.0 | | 5.7 | 0.67 |
| 56-23-5 | Carbon tetrachloride | 24.0 | | 5.7 | 0.74 |
| 108-90-7 | Chlorobenzene | 20.7 | | 5.7 | 0.60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 MS | Lab Sample ID: 240-57597-5 MS

Matrix: Solid | Lab File ID: 146341.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.17(g) | Date Analyzed: 11/10/2015 16:30

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 20.2 | | 5.7 | 0.42 |
| 75-00-3 | Chloroethane | 21.5 | | 5.7 | 0.50 |
| 67-66-3 | Chloroform | 22.9 | | 5.7 | 0.41 |
| 74-87-3 | Chloromethane | 15.9 | | 5.7 | 0.90 |
| 156-59-2 | cis-1,2-Dichloroethene | 22.1 | | 5.7 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 19.1 | | 5.7 | 0.81 |
| 74-95-3 | Dibromomethane | 22.0 | | 5.7 | 0.48 |
| 74-97-5 | Bromochloromethane | 23.3 | | 5.7 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 18.3 | | 5.7 | 0.32 |
| 75-27-4 | Bromodichloromethane | 20.0 | | 5.7 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 164 | J | 230 | 6.5 |
| 100-41-4 | Ethylbenzene | 20.5 | | 5.7 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 12.5 | | 5.7 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.67 | U | 11 | 0.67 |
| 98-82-8 | Isopropylbenzene | 19.5 | | 5.7 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 20.9 | | 5.7 | 0.49 |
| 75-09-2 | Methylene Chloride | 18.8 | | 5.7 | 0.84 |
| 179601-23-1 | m-Xylene & p-Xylene | 19.3 | | 11 | 0.48 |
| 91-20-3 | Naphthalene | 17.8 | | 5.7 | 0.40 |
| 104-51-8 | n-Butylbenzene | 16.6 | | 5.7 | 0.67 |
| 103-65-1 | N-Propylbenzene | 19.7 | | 5.7 | 0.68 |
| 95-47-6 | o-Xylene | 20.9 | | 5.7 | 0.65 |
| 135-98-8 | sec-Butylbenzene | 18.2 | | 5.7 | 1.0 |
| 100-42-5 | Styrene | 17.9 | | 5.7 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.49 | U | 5.7 | 0.49 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.7 | 0.47 |
| 127-18-4 | Tetrachloroethene | 22.3 | | 5.7 | 0.91 |
| 108-88-3 | Toluene | 22.0 | | 5.7 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 22.3 | | 5.7 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 21.1 | | 5.7 | 0.33 |
| 79-01-6 | Trichloroethene | 23.2 | | 5.7 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 23.0 | | 5.7 | 0.33 |
| 108-05-4 | Vinyl acetate | 14.4 | | 11 | 0.75 |
| 75-01-4 | Vinyl chloride | 17.4 | | 5.7 | 0.34 |
| 1330-20-7 | Xylenes, Total | 40.2 | | 11 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 21.3 | | 5.7 | 0.76 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 MS | Lab Sample ID: 240-57597-5 MS

Matrix: Solid | Lab File ID: 146341.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.17(g) | Date Analyzed: 11/10/2015 16:30

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 24.2 | | 5.7 | 0.60 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 99 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 97 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 95 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 92 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: 146366.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.91(g) | Date Analyzed: 11/11/2015 13:31

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 20.2 | | 4.9 | 0.49 |
| 71-55-6 | 1,1,1-Trichloroethane | 22.0 | | 4.9 | 0.64 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 19.2 | | 4.9 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 21.4 | | 4.9 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 18.4 | | 4.9 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 18.3 | | 4.9 | 0.79 |
| 563-58-6 | 1,1-Dichloropropene | 20.4 | | 4.9 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 19.1 | | 4.9 | 0.49 |
| 96-18-4 | 1,2,3-Trichloropropane | 22.1 | | 4.9 | 0.77 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 4.9 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 19.5 | | 4.9 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 19.4 | | 4.9 | 0.57 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 17.8 | | 9.9 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 18.5 | | 4.9 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 19.3 | | 4.9 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 17.0 | | 4.9 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 19.2 | | 4.9 | 0.68 |
| 142-28-9 | 1,3-Dichloropropane | 18.5 | | 4.9 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 17.9 | | 4.9 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 21.6 | | 4.9 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 30.1 | | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 12.6 | J | 49 | 0.63 |
| 95-49-8 | 2-Chlorotoluene | 19.9 | | 4.9 | 0.30 |
| 591-78-6 | 2-Hexanone | 31.1 | | 20 | 0.67 |
| 106-43-4 | 4-Chlorotoluene | 19.3 | | 4.9 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 20.1 | | 4.9 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 32.3 | | 20 | 1.2 |
| 67-64-1 | Acetone | 30.0 | | 20 | 5.6 |
| 71-43-2 | Benzene | 18.5 | | 4.9 | 0.89 |
| 108-86-1 | Bromobenzene | 19.9 | | 4.9 | 0.33 |
| 75-25-2 | Bromoform | 16.1 | | 4.9 | 0.24 |
| 74-83-9 | Bromomethane | 19.1 | | 4.9 | 0.42 |
| 75-15-0 | Carbon disulfide | 13.8 | | 4.9 | 0.58 |
| 56-23-5 | Carbon tetrachloride | 21.4 | | 4.9 | 0.64 |
| 108-90-7 | Chlorobenzene | 19.3 | | 4.9 | 0.52 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: 146366.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.91(g) | Date Analyzed: 11/11/2015 13:31

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 17.1 | | 4.9 | 0.37 |
| 75-00-3 | Chloroethane | 19.4 | | 4.9 | 0.44 |
| 67-66-3 | Chloroform | 20.5 | | 4.9 | 0.36 |
| 74-87-3 | Chloromethane | 15.5 | | 4.9 | 0.78 |
| 156-59-2 | cis-1,2-Dichloroethene | 19.6 | | 4.9 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 16.0 | | 4.9 | 0.70 |
| 74-95-3 | Dibromomethane | 21.4 | | 4.9 | 0.42 |
| 74-97-5 | Bromochloromethane | 20.6 | | 4.9 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 19.9 | | 4.9 | 0.28 |
| 75-27-4 | Bromodichloromethane | 17.2 | | 4.9 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 163 | J | 200 | 5.6 |
| 100-41-4 | Ethylbenzene | 19.0 | | 4.9 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 20.9 | | 4.9 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.58 | U | 9.9 | 0.58 |
| 98-82-8 | Isopropylbenzene | 20.0 | | 4.9 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 17.5 | | 4.9 | 0.43 |
| 75-09-2 | Methylene Chloride | 16.2 | | 4.9 | 0.73 |
| 179601-23-1 | m-Xylene & p-Xylene | 19.4 | | 9.9 | 0.42 |
| 91-20-3 | Naphthalene | 16.9 | | 4.9 | 0.35 |
| 104-51-8 | n-Butylbenzene | 19.3 | | 4.9 | 0.58 |
| 103-65-1 | N-Propylbenzene | 19.4 | | 4.9 | 0.59 |
| 95-47-6 | o-Xylene | 19.5 | | 4.9 | 0.56 |
| 135-98-8 | sec-Butylbenzene | 19.5 | | 4.9 | 0.90 |
| 100-42-5 | Styrene | 17.8 | | 4.9 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 4.9 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 4.9 | 0.41 |
| 127-18-4 | Tetrachloroethene | 21.7 | | 4.9 | 0.79 |
| 108-88-3 | Toluene | 20.2 | | 4.9 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 19.7 | | 4.9 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 17.1 | | 4.9 | 0.29 |
| 79-01-6 | Trichloroethene | 20.3 | | 4.9 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 21.2 | | 4.9 | 0.29 |
| 108-05-4 | Vinyl acetate | 17.6 | | 9.9 | 0.65 |
| 75-01-4 | Vinyl chloride | 16.9 | | 4.9 | 0.30 |
| 1330-20-7 | Xylenes, Total | 38.9 | | 9.9 | 0.53 |
| 98-06-6 | tert-Butylbenzene | 20.0 | | 4.9 | 0.66 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: 146366.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.91(g) | Date Analyzed: 11/11/2015 13:31

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 19.8 | | 4.9 | 0.52 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 97 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 97 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 88 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 97 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57564-B-1 MS

Matrix: Water | Lab File ID: UXC4417.D

Analysis Method: 8260C | Date Collected: 11/05/2015 12:10

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 250

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 71-55-6 | 1,1,1-Trichloroethane | 2570 | | 250 | 110 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 2130 | | 250 | 55 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 1850 | | 250 | 110 |
| 75-34-3 | 1,1-Dichloroethane | 2270 | | 250 | 75 |
| 75-35-4 | 1,1-Dichloroethene | 2020 | | 250 | 110 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 2090 | | 250 | 80 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2310 | | 500 | 210 |
| 95-50-1 | 1,2-Dichlorobenzene | 2300 | | 250 | 63 |
| 107-06-2 | 1,2-Dichloroethane | 2580 | | 250 | 58 |
| 78-87-5 | 1,2-Dichloropropane | 2170 | | 250 | 63 |
| 541-73-1 | 1,3-Dichlorobenzene | 2250 | | 250 | 48 |
| 106-46-7 | 1,4-Dichlorobenzene | 2290 | | 250 | 68 |
| 78-93-3 | 2-Butanone (MEK) | 12400 | | 2500 | 130 |
| 591-78-6 | 2-Hexanone | 4460 | | 2500 | 120 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5380 | | 2500 | 250 |
| 67-64-1 | Acetone | 18700 | | 2500 | 240 |
| 71-43-2 | Benzene | 2220 | | 250 | 88 |
| 75-25-2 | Bromoform | 2500 | | 250 | 140 |
| 74-83-9 | Bromomethane | 2390 | | 250 | 110 |
| 75-15-0 | Carbon disulfide | 1870 | | 250 | 95 |
| 56-23-5 | Carbon tetrachloride | 2620 | | 250 | 110 |
| 108-90-7 | Chlorobenzene | 2390 | | 250 | 63 |
| 124-48-1 | Dibromochloromethane | 2460 | | 250 | 110 |
| 75-00-3 | Chloroethane | 2200 | | 250 | 80 |
| 67-66-3 | Chloroform | 2560 | | 250 | 63 |
| 74-87-3 | Chloromethane | 1980 | | 250 | 110 |
| 156-59-2 | cis-1,2-Dichloroethene | 2370 | | 250 | 65 |
| 10061-01-5 | cis-1,3-Dichloropropene | 2250 | | 250 | 120 |
| 75-71-8 | Dichlorodifluoromethane | 1870 | | 250 | 80 |
| 75-27-4 | Bromodichloromethane | 2650 | | 250 | 73 |
| 100-41-4 | Ethylbenzene | 2440 | | 250 | 63 |
| 98-82-8 | Isopropylbenzene | 2380 | | 250 | 88 |
| 1634-04-4 | Methyl tert-butyl ether | 2520 | | 250 | 50 |
| 75-09-2 | Methylene Chloride | 2350 | | 250 | 83 |
| 179601-23-1 | m-Xylene & p-Xylene | 2540 | | 500 | 60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57564-B-1 MS

Matrix: Water | Lab File ID: UXC4417.D

Analysis Method: 8260C | Date Collected: 11/05/2015 12:10

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 250

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 95-47-6 | o-Xylene | 2500 | | 250 | 63 |
| 100-42-5 | Styrene | 2320 | | 250 | 110 |
| 127-18-4 | Tetrachloroethene | 2450 | | 250 | 78 |
| 108-88-3 | Toluene | 2380 | | 250 | 58 |
| 156-60-5 | trans-1,2-Dichloroethene | 2340 | | 250 | 75 |
| 10061-02-6 | trans-1,3-Dichloropropene | 2330 | | 250 | 140 |
| 79-01-6 | Trichloroethene | 2580 | | 250 | 55 |
| 75-69-4 | Trichlorofluoromethane | 2020 | | 250 | 120 |
| 75-01-4 | Vinyl chloride | 2040 | | 250 | 73 |
| 1330-20-7 | Xylenes, Total | 5040 | | 500 | 130 |
| 106-93-4 | 1,2-Dibromoethane | 2430 | | 250 | 80 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 88 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 99 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 103 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 97 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 MSD | Lab Sample ID: 240-57597-5 MSD

Matrix: Solid | Lab File ID: 146342.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.12(g) | Date Analyzed: 11/10/2015 16:52

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 24.0 | | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 25.7 | | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 28.0 | | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 25.6 | | 5.8 | 0.48 |
| 75-34-3 | 1,1-Dichloroethane | 22.5 | | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 22.0 | | 5.8 | 0.92 |
| 563-58-6 | 1,1-Dichloropropene | 25.1 | | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 17.9 | | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 28.9 | | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 1.64 | J | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 19.9 | | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 25.2 | | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 26.9 | | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 22.2 | | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 24.1 | | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 21.9 | | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 22.6 | | 5.8 | 0.79 |
| 142-28-9 | 1,3-Dichloropropane | 23.2 | | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 22.3 | | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 24.5 | | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 39.3 | | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 17.2 | J | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 25.6 | | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 52.6 | | 23 | 0.78 |
| 106-43-4 | 4-Chlorotoluene | 24.9 | | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 26.0 | | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 140 | | 23 | 1.4 |
| 67-64-1 | Acetone | 53.3 | | 23 | 6.6 |
| 71-43-2 | Benzene | 23.0 | | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 26.9 | | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 19.9 | | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 21.9 | | 5.8 | 0.48 |
| 75-15-0 | Carbon disulfide | 17.6 | | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 26.5 | | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 23.7 | | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 MSD | Lab Sample ID: 240-57597-5 MSD

Matrix: Solid | Lab File ID: 146342.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.12(g) | Date Analyzed: 11/10/2015 16:52

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 21.2 | | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 22.3 | | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 25.1 | | 5.8 | 0.41 |
| 74-87-3 | Chloromethane | 17.1 | | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 24.0 | | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 21.0 | | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 24.2 | | 5.8 | 0.48 |
| 74-97-5 | Bromochloromethane | 26.0 | | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 18.5 | | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 22.9 | | 5.8 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 189 | J | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 24.4 | | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 22.3 | | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 25.1 | | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 22.3 | | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 20.2 | | 5.8 | 0.85 |
| 179601-23-1 | m-Xylene & p-Xylene | 24.0 | | 12 | 0.48 |
| 91-20-3 | Naphthalene | 21.0 | | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 25.0 | | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 25.5 | | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 25.1 | | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 25.1 | | 5.8 | 1.0 |
| 100-42-5 | Styrene | 21.9 | | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 24.7 | | 5.8 | 0.92 |
| 108-88-3 | Toluene | 24.4 | | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 24.7 | | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 22.9 | | 5.8 | 0.33 |
| 79-01-6 | Trichloroethene | 25.3 | | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 26.2 | | 5.8 | 0.33 |
| 108-05-4 | Vinyl acetate | 17.8 | | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 18.6 | | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 49.1 | | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 26.7 | | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 MSD | Lab Sample ID: 240-57597-5 MSD

Matrix: Solid | Lab File ID: 146342.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.12(g) | Date Analyzed: 11/10/2015 16:52

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 25.0 | | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 101 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 96 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 98 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD | Lab Sample ID: 240-57597-15 MSD

Matrix: Solid | Lab File ID: 146367.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.17(g) | Date Analyzed: 11/11/2015 13:52

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 22.0 | | 5.7 | 0.57 |
| 71-55-6 | 1,1,1-Trichloroethane | 23.8 | | 5.7 | 0.74 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 21.3 | | 5.7 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 23.5 | | 5.7 | 0.47 |
| 75-34-3 | 1,1-Dichloroethane | 20.4 | | 5.7 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 20.3 | | 5.7 | 0.90 |
| 563-58-6 | 1,1-Dichloropropene | 22.2 | | 5.7 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 19.9 | | 5.7 | 0.57 |
| 96-18-4 | 1,2,3-Trichloropropane | 20.5 | | 5.7 | 0.88 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.52 | U | 5.7 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 20.8 | | 5.7 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 20.8 | | 5.7 | 0.66 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 20.5 | | 11 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 19.9 | | 5.7 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 21.1 | | 5.7 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 19.2 | | 5.7 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 19.0 | | 5.7 | 0.78 |
| 142-28-9 | 1,3-Dichloropropane | 20.2 | | 5.7 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 19.1 | | 5.7 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 24.1 | | 5.7 | 0.50 |
| 78-93-3 | 2-Butanone (MEK) | 31.6 | | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 14.8 | J | 57 | 0.72 |
| 95-49-8 | 2-Chlorotoluene | 21.8 | | 5.7 | 0.34 |
| 591-78-6 | 2-Hexanone | 33.3 | | 23 | 0.77 |
| 106-43-4 | 4-Chlorotoluene | 20.1 | | 5.7 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 21.4 | | 5.7 | 0.46 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 33.2 | | 23 | 1.4 |
| 67-64-1 | Acetone | 29.1 | | 23 | 6.4 |
| 71-43-2 | Benzene | 20.6 | | 5.7 | 1.0 |
| 108-86-1 | Bromobenzene | 22.2 | | 5.7 | 0.37 |
| 75-25-2 | Bromoform | 17.5 | | 5.7 | 0.27 |
| 74-83-9 | Bromomethane | 20.8 | | 5.7 | 0.47 |
| 75-15-0 | Carbon disulfide | 15.4 | | 5.7 | 0.67 |
| 56-23-5 | Carbon tetrachloride | 23.7 | | 5.7 | 0.74 |
| 108-90-7 | Chlorobenzene | 20.8 | | 5.7 | 0.60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD | Lab Sample ID: 240-57597-15 MSD

Matrix: Solid | Lab File ID: 146367.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.17(g) | Date Analyzed: 11/11/2015 13:52

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 19.4 | | 5.7 | 0.42 |
| 75-00-3 | Chloroethane | 20.8 | | 5.7 | 0.50 |
| 67-66-3 | Chloroform | 22.6 | | 5.7 | 0.41 |
| 74-87-3 | Chloromethane | 17.2 | | 5.7 | 0.89 |
| 156-59-2 | cis-1,2-Dichloroethene | 21.5 | | 5.7 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 18.1 | | 5.7 | 0.80 |
| 74-95-3 | Dibromomethane | 22.8 | | 5.7 | 0.47 |
| 74-97-5 | Bromochloromethane | 22.2 | | 5.7 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 21.8 | | 5.7 | 0.32 |
| 75-27-4 | Bromodichloromethane | 20.4 | | 5.7 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 158 | J | 230 | 6.5 |
| 100-41-4 | Ethylbenzene | 20.9 | | 5.7 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 22.4 | | 5.7 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.67 | U | 11 | 0.67 |
| 98-82-8 | Isopropylbenzene | 21.8 | | 5.7 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 19.5 | | 5.7 | 0.49 |
| 75-09-2 | Methylene Chloride | 19.2 | | 5.7 | 0.84 |
| 179601-23-1 | m-Xylene & p-Xylene | 20.2 | | 11 | 0.47 |
| 91-20-3 | Naphthalene | 18.8 | | 5.7 | 0.40 |
| 104-51-8 | n-Butylbenzene | 20.7 | | 5.7 | 0.67 |
| 103-65-1 | N-Propylbenzene | 20.7 | | 5.7 | 0.68 |
| 95-47-6 | o-Xylene | 21.5 | | 5.7 | 0.64 |
| 135-98-8 | sec-Butylbenzene | 20.6 | | 5.7 | 1.0 |
| 100-42-5 | Styrene | 19.5 | | 5.7 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.49 | U | 5.7 | 0.49 |
| 637-92-3 | Ethyl t-butyl ether | 0.46 | U | 5.7 | 0.46 |
| 127-18-4 | Tetrachloroethene | 21.7 | | 5.7 | 0.90 |
| 108-88-3 | Toluene | 21.7 | | 5.7 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 21.8 | | 5.7 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 19.7 | | 5.7 | 0.33 |
| 79-01-6 | Trichloroethene | 22.5 | | 5.7 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 25.1 | | 5.7 | 0.33 |
| 108-05-4 | Vinyl acetate | 17.9 | | 11 | 0.75 |
| 75-01-4 | Vinyl chloride | 18.5 | | 5.7 | 0.34 |
| 1330-20-7 | Xylenes, Total | 41.7 | | 11 | 0.61 |
| 98-06-6 | tert-Butylbenzene | 21.5 | | 5.7 | 0.76 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD | Lab Sample ID: 240-57597-15 MSD

Matrix: Solid | Lab File ID: 146367.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.17(g) | Date Analyzed: 11/11/2015 13:52

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 21.9 | | 5.7 | 0.60 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 96 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 99 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 86 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 95 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57564-B-1 MSD

Matrix: Water | Lab File ID: UXC4418.D

Analysis Method: 8260C | Date Collected: 11/05/2015 12:10

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 20:27

Soil Aliquot Vol: | Dilution Factor: 250

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 71-55-6 | 1,1,1-Trichloroethane | 2630 | | 250 | 110 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 2220 | | 250 | 55 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2160 | | 250 | 110 |
| 75-34-3 | 1,1-Dichloroethane | 2310 | | 250 | 75 |
| 75-35-4 | 1,1-Dichloroethene | 2090 | | 250 | 110 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 2190 | | 250 | 80 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2510 | | 500 | 210 |
| 95-50-1 | 1,2-Dichlorobenzene | 2370 | | 250 | 63 |
| 107-06-2 | 1,2-Dichloroethane | 2590 | | 250 | 58 |
| 78-87-5 | 1,2-Dichloropropane | 2330 | | 250 | 63 |
| 541-73-1 | 1,3-Dichlorobenzene | 2270 | | 250 | 48 |
| 106-46-7 | 1,4-Dichlorobenzene | 2330 | | 250 | 68 |
| 78-93-3 | 2-Butanone (MEK) | 12800 | | 2500 | 130 |
| 591-78-6 | 2-Hexanone | 4330 | | 2500 | 120 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 5460 | | 2500 | 250 |
| 67-64-1 | Acetone | 18900 | | 2500 | 240 |
| 71-43-2 | Benzene | 2250 | | 250 | 88 |
| 75-25-2 | Bromoform | 2490 | | 250 | 140 |
| 74-83-9 | Bromomethane | 2400 | | 250 | 110 |
| 75-15-0 | Carbon disulfide | 1920 | | 250 | 95 |
| 56-23-5 | Carbon tetrachloride | 2790 | | 250 | 110 |
| 108-90-7 | Chlorobenzene | 2400 | | 250 | 63 |
| 124-48-1 | Dibromochloromethane | 2450 | | 250 | 110 |
| 75-00-3 | Chloroethane | 2180 | | 250 | 80 |
| 67-66-3 | Chloroform | 2580 | | 250 | 63 |
| 74-87-3 | Chloromethane | 2060 | | 250 | 110 |
| 156-59-2 | cis-1,2-Dichloroethene | 2420 | | 250 | 65 |
| 10061-01-5 | cis-1,3-Dichloropropene | 2330 | | 250 | 120 |
| 75-71-8 | Dichlorodifluoromethane | 2230 | | 250 | 80 |
| 75-27-4 | Bromodichloromethane | 2760 | | 250 | 73 |
| 100-41-4 | Ethylbenzene | 2370 | | 250 | 63 |
| 98-82-8 | Isopropylbenzene | 2400 | | 250 | 88 |
| 1634-04-4 | Methyl tert-butyl ether | 2570 | | 250 | 50 |
| 75-09-2 | Methylene Chloride | 2330 | | 250 | 83 |
| 179601-23-1 | m-Xylene & p-Xylene | 2510 | | 500 | 60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57564-B-1 MSD

Matrix: Water | Lab File ID: UXC4418.D

Analysis Method: 8260C | Date Collected: 11/05/2015 12:10

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 20:27

Soil Aliquot Vol: | Dilution Factor: 250

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 95-47-6 | o-Xylene | 2490 | | 250 | 63 |
| 100-42-5 | Styrene | 2340 | | 250 | 110 |
| 127-18-4 | Tetrachloroethene | 2490 | | 250 | 78 |
| 108-88-3 | Toluene | 2430 | | 250 | 58 |
| 156-60-5 | trans-1,2-Dichloroethene | 2400 | | 250 | 75 |
| 10061-02-6 | trans-1,3-Dichloropropene | 2310 | | 250 | 140 |
| 79-01-6 | Trichloroethene | 2730 | | 250 | 55 |
| 75-69-4 | Trichlorofluoromethane | 2240 | | 250 | 120 |
| 75-01-4 | Vinyl chloride | 2050 | | 250 | 73 |
| 1330-20-7 | Xylenes, Total | 5000 | | 500 | 130 |
| 106-93-4 | 1,2-Dibromoethane | 2470 | | 250 | 80 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 90 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 102 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 98 | | 78-125 |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A3UX14 Start Date: 08/13/2015 13:54

Analysis Batch Number: 193150 End Date: 08/13/2015 23:12

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-193150/28 | | 08/13/2015 13:54 | 1 | BFB14045.D | DB-624 0.18(mm) |
| IC 240-193150/3 | | 08/13/2015 14:36 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/4 | | 08/13/2015 14:58 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/5 | | 08/13/2015 15:20 | 1 | | DB-624 0.18(mm) |
| ICIS 240-193150/6 | | 08/13/2015 15:41 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/7 | | 08/13/2015 16:03 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/8 | | 08/13/2015 16:24 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/9 | | 08/13/2015 16:45 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/10 | | 08/13/2015 17:07 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/11 | | 08/13/2015 17:28 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/12 | | 08/13/2015 17:50 | 1 | 144576.D | DB-624 0.18(mm) |
| IC 240-193150/13 | | 08/13/2015 18:11 | 1 | 144577.D | DB-624 0.18(mm) |
| IC 240-193150/14 | | 08/13/2015 18:33 | 1 | 144578.D | DB-624 0.18(mm) |
| IC 240-193150/15 | | 08/13/2015 18:54 | 1 | 144579.D | DB-624 0.18(mm) |
| IC 240-193150/16 | | 08/13/2015 19:16 | 1 | 144580.D | DB-624 0.18(mm) |
| IC 240-193150/17 | | 08/13/2015 19:37 | 1 | 144581.D | DB-624 0.18(mm) |
| IC 240-193150/18 | | 08/13/2015 19:59 | 1 | 144582.D | DB-624 0.18(mm) |
| IC 240-193150/19 | | 08/13/2015 20:20 | 1 | 144583.D | DB-624 0.18(mm) |
| IC 240-193150/20 | | 08/13/2015 20:42 | 1 | 144584.D | DB-624 0.18(mm) |
| ICV 240-193150/21 | | 08/13/2015 21:03 | 1 | | DB-624 0.18(mm) |
| ICV 240-193150/22 | | 08/13/2015 21:25 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 21:46 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 22:08 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 22:29 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 22:50 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 23:12 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A3UX14 Start Date: 09/02/2015 13:49

Analysis Batch Number: 196026 End Date: 09/02/2015 19:33

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-196026/1 | | 09/02/2015 13:49 | 1 | BFB14070.D | DB-624 0.18(mm) |
| IC 240-196026/3 | | 09/02/2015 14:32 | 1 | 144935.D | DB-624 0.18(mm) |
| IC 240-196026/4 | | 09/02/2015 14:54 | 1 | 144936.D | DB-624 0.18(mm) |
| IC 240-196026/5 | | 09/02/2015 15:15 | 1 | 144937.D | DB-624 0.18(mm) |
| ICIS 240-196026/6 | | 09/02/2015 15:37 | 1 | 144938.D | DB-624 0.18(mm) |
| IC 240-196026/7 | | 09/02/2015 15:58 | 1 | 144939.D | DB-624 0.18(mm) |
| IC 240-196026/8 | | 09/02/2015 16:20 | 1 | 144940.D | DB-624 0.18(mm) |
| IC 240-196026/9 | | 09/02/2015 16:41 | 1 | 144941.D | DB-624 0.18(mm) |
| IC 240-196026/10 | | 09/02/2015 17:03 | 1 | 144942.D | DB-624 0.18(mm) |
| IC 240-196026/11 | | 09/02/2015 17:24 | 1 | 144943.D | DB-624 0.18(mm) |
| CCV 240-196026/12 | | 09/02/2015 17:46 | 1 | | DB-624 0.18(mm) |
| ICV 240-196026/13 | | 09/02/2015 18:07 | 1 | 144945.D | DB-624 0.18(mm) |
| ZZZZZ | | 09/02/2015 18:29 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 09/02/2015 18:50 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 09/02/2015 19:12 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 09/02/2015 19:33 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A3UX14 Start Date: 11/10/2015 10:04

Analysis Batch Number: 206051 End Date: 11/10/2015 20:05

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-206051/1 | | 11/10/2015 10:04 | 1 | BFB14128.D | DB-624 0.18(mm) |
| CCVIS 240-206051/3 | | 11/10/2015 10:47 | 1 | 146325.D | DB-624 0.18(mm) |
| CCV 240-206051/4 | | 11/10/2015 11:08 | 1 | 146326.D | DB-624 0.18(mm) |
| LCS 240-206051/5 | | 11/10/2015 11:30 | 1 | 146327.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/10/2015 11:51 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/10/2015 12:13 | 1 | | DB-624 0.18(mm) |
| MB 240-206089/1-A | | 11/10/2015 12:56 | 1 | 146331.D | DB-624 0.18(mm) |
| 240-57597-1 | F-VB-UST-1-14 | 11/10/2015 14:43 | 1 | 146336.D | DB-624 0.18(mm) |
| 240-57597-2 | F-VB-UST-2-14 | 11/10/2015 15:05 | 1 | 146337.D | DB-624 0.18(mm) |
| 240-57597-3 | F-VB-UST-3-14 | 11/10/2015 15:26 | 1 | 146338.D | DB-624 0.18(mm) |
| 240-57597-4 | F-VB-UST-4-14 | 11/10/2015 15:48 | 1 | 146339.D | DB-624 0.18(mm) |
| 240-57597-5 | F-VB-UST-5-14 | 11/10/2015 16:09 | 1 | 146340.D | DB-624 0.18(mm) |
| 240-57597-5 MS | F-VB-UST-5-14 MS | 11/10/2015 16:30 | 1 | 146341.D | DB-624 0.18(mm) |
| 240-57597-5 MSD | F-VB-UST-5-14 MSD | 11/10/2015 16:52 | 1 | 146342.D | DB-624 0.18(mm) |
| 240-57597-6 | F-VB-UST-6-14 | 11/10/2015 17:13 | 1 | 146343.D | DB-624 0.18(mm) |
| 240-57597-7 | F-VB-UST-7-14 | 11/10/2015 17:35 | 1 | 146344.D | DB-624 0.18(mm) |
| 240-57597-8 | F-VB-UST-8-14 | 11/10/2015 17:56 | 1 | 146345.D | DB-624 0.18(mm) |
| 240-57597-9 | F-VB-UST-9-14 | 11/10/2015 18:18 | 1 | 146346.D | DB-624 0.18(mm) |
| 240-57597-10 | F-VB-UST-10-14 | 11/10/2015 18:39 | 1 | 146347.D | DB-624 0.18(mm) |
| 240-57597-11 | F-VB-UST-11-14 | 11/10/2015 19:01 | 1 | 146348.D | DB-624 0.18(mm) |
| 240-57597-12 | F-VB-UST-12-14 | 11/10/2015 19:22 | 1 | 146349.D | DB-624 0.18(mm) |
| 240-57597-13 | F-VB-UST-13-14 | 11/10/2015 19:44 | 1 | 146350.D | DB-624 0.18(mm) |
| 240-57597-14 | F-VB-UST-14-14 | 11/10/2015 20:05 | 1 | 146351.D | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A3UX14 Start Date: 11/11/2015 09:51

Analysis Batch Number: 206267 End Date: 11/11/2015 14:14

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-206267/1 | | 11/11/2015 09:51 | 1 | BFB14129.D | DB-624 0.18(mm) |
| CCVIS 240-206267/3 | | 11/11/2015 10:39 | 1 | 146358.D | DB-624 0.18(mm) |
| CCV 240-206267/4 | | 11/11/2015 11:00 | 1 | 146359.D | DB-624 0.18(mm) |
| LCS 240-206267/5 | | 11/11/2015 11:22 | 1 | 146360.D | DB-624 0.18(mm) |
| MB 240-206267/6 | | 11/11/2015 11:43 | 1 | 146361.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/11/2015 12:05 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/11/2015 12:26 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/11/2015 12:48 | 1 | | DB-624 0.18(mm) |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 13:09 | 1 | 146365.D | DB-624 0.18(mm) |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 13:31 | 1 | 146366.D | DB-624 0.18(mm) |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 13:52 | 1 | 146367.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/11/2015 14:14 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A3UX15 Start Date: 10/28/2015 12:37

Analysis Batch Number: 204067 End Date: 10/28/2015 18:00

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-204067/1 | | 10/28/2015 12:37 | 1 | BFB460.D | DB-624 0.18(mm) |
| STD8260 240-204067/2 IC | | 10/28/2015 13:05 | 1 | UXC3824A.D | DB-624 0.18(mm) |
| STD8260 240-204067/3 IC | | 10/28/2015 13:28 | 1 | UXC3825A.D | DB-624 0.18(mm) |
| STD8260 240-204067/4 ICIS | | 10/28/2015 13:51 | 1 | UXC3826A.D | DB-624 0.18(mm) |
| STD8260 240-204067/5 IC | | 10/28/2015 14:13 | 1 | UXC3827A.D | DB-624 0.18(mm) |
| STD8260 240-204067/6 IC | | 10/28/2015 14:36 | 1 | UXC3828A.D | DB-624 0.18(mm) |
| STD8260 240-204067/7 IC | | 10/28/2015 14:59 | 1 | UXC3829A.D | DB-624 0.18(mm) |
| ICV 240-204067/14 | | 10/28/2015 15:21 | 1 | UXC3830.D | DB-624 0.18(mm) |
| STD 240-204067/8 IC | | 10/28/2015 15:44 | 1 | UXC3831.D | DB-624 0.18(mm) |
| STD 240-204067/9 IC | | 10/28/2015 16:06 | 1 | UXC3832.D | DB-624 0.18(mm) |
| STD 240-204067/10 IC | | 10/28/2015 16:29 | 1 | UXC3833.D | DB-624 0.18(mm) |
| STD 240-204067/11 IC | | 10/28/2015 16:52 | 1 | UXC3834.D | DB-624 0.18(mm) |
| STD 240-204067/12 IC | | 10/28/2015 17:14 | 1 | UXC3835.D | DB-624 0.18(mm) |
| STD 240-204067/13 IC | | 10/28/2015 17:37 | 1 | UXC3836.D | DB-624 0.18(mm) |
| ICV 240-204067/15 | | 10/28/2015 18:00 | 1 | UXC3837.D | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Instrument ID: A3UX15  Start Date: 11/12/2015 10:14

Analysis Batch Number: 206461  End Date: 11/12/2015 20:27

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-206461/1 | | 11/12/2015 10:14 | 1 | BFB472.D | DB-624 0.18(mm) |
| CCVIS 240-206461/2 | | 11/12/2015 10:41 | 1 | UXC4392.D | DB-624 0.18(mm) |
| LCS 240-206461/4 | | 11/12/2015 11:03 | 1 | UXC4393.D | DB-624 0.18(mm) |
| CCV 240-206461/3 | | 11/12/2015 11:26 | 1 | UXC4394.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/12/2015 11:48 | 1 | | DB-624 0.18(mm) |
| MB 240-206461/6 | | 11/12/2015 12:11 | 1 | UXC4396.D | DB-624 0.18(mm) |
| 240-57597-16 | TB-110515 | 11/12/2015 12:33 | 1 | UXC4397.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/12/2015 12:56 | 250 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/12/2015 13:19 | 250 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/12/2015 13:41 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/12/2015 14:04 | 1 | | DB-624 0.18(mm) |
| 240-57564-B-1 MS | | 11/12/2015 20:05 | 250 | UXC4417.D | DB-624 0.18(mm) |
| 240-57564-B-1 MSD | | 11/12/2015 20:27 | 250 | UXC4418.D | DB-624 0.18(mm) |

# GC/MS VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 206089  Batch Start Date: 11/10/15 11:24  Batch Analyst: Macenczak, Steven

Batch Method: 5030C  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| MB 240-206089/1 | | 5030C, 8260C | | 5.0 g | 5 mL | | | | |
| 240-57597-B-1 | F-VB-UST-1-14 | 5030C, 8260C | T | 5.89 g | 5 mL | | | | |
| 240-57597-B-2 | F-VB-UST-2-14 | 5030C, 8260C | T | 5.36 g | 5 mL | | | | |
| 240-57597-B-3 | F-VB-UST-3-14 | 5030C, 8260C | T | 5.23 g | 5 mL | | | | |
| 240-57597-B-4 | F-VB-UST-4-14 | 5030C, 8260C | T | 5.43 g | 5 mL | | | | |
| 240-57597-B-5 | F-VB-UST-5-14 | 5030C, 8260C | T | 5.09 g | 5 mL | | | | |
| 240-57597-B-5 MS | F-VB-UST-5-14 | 5030C, 8260C | T | 5.17 g | 5 mL | | | | |
| 240-57597-B-5 MSD | F-VB-UST-5-14 | 5030C, 8260C | T | 5.12 g | 5 mL | | | | |
| 240-57597-B-6 | F-VB-UST-6-14 | 5030C, 8260C | T | 5.18 g | 5 mL | | | | |
| 240-57597-B-7 | F-VB-UST-7-14 | 5030C, 8260C | T | 5.31 g | 5 mL | | | | |
| 240-57597-B-8 | F-VB-UST-8-14 | 5030C, 8260C | T | 5.29 g | 5 mL | | | | |
| 240-57597-B-9 | F-VB-UST-9-14 | 5030C, 8260C | T | 5.35 g | 5 mL | | | | |
| 240-57597-B-10 | F-VB-UST-10-14 | 5030C, 8260C | T | 5.14 g | 5 mL | | | | |
| 240-57597-B-11 | F-VB-UST-11-14 | 5030C, 8260C | T | 5.03 g | 5 mL | | | | |
| 240-57597-B-12 | F-VB-UST-12-14 | 5030C, 8260C | T | 5.37 g | 5 mL | | | | |
| 240-57597-B-13 | F-VB-UST-13-14 | 5030C, 8260C | T | 5.67 g | 5 mL | | | | |
| 240-57597-B-14 | F-VB-UST-14-14 | 5030C, 8260C | T | 5.55 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| | |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC/MS VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 206305  Batch Start Date: 11/11/15 11:03  Batch Analyst: Macenczak, Steven

Batch Method: 5030C  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-B-15 | F-VB-UST-DUP | 5030C, 8260C | T | 5.55 g | 5 mL | | | | |
| 240-57597-B-15 MS | F-VB-UST-DUP | 5030C, 8260C | T | 5.91 g | 5 mL | | | | |
| 240-57597-B-15 MSD | F-VB-UST-DUP | 5030C, 8260C | T | 5.17 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| | |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8270D

## Semivolatile Organic Compounds (GC/MS) by Method 8270D

FORM II
GC/MS SEMI VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): DB-624 ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | 2FP # | PHL # | NBZ # | FBP # | TBP # | TPH # |
|---|---|---|---|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 64 | 65 | 62 | 67 | 44 | 76 |
| F-VB-UST-2-14 | 240-57597-2 | 61 | 61 | 58 | 63 | 42 | 78 |
| F-VB-UST-3-14 | 240-57597-3 | 61 | 60 | 58 | 63 | 41 | 73 |
| F-VB-UST-4-14 | 240-57597-4 | 72 | 72 | 69 | 72 | 42 | 85 |
| F-VB-UST-5-14 | 240-57597-5 | 68 | 67 | 64 | 69 | 25 | 77 |
| F-VB-UST-6-14 | 240-57597-6 | 62 | 60 | 62 | 68 | 33 | 79 |
| F-VB-UST-7-14 | 240-57597-7 | 67 | 64 | 62 | 70 | 38 | 80 |
| F-VB-UST-8-14 | 240-57597-8 | 56 | 54 | 52 | 56 | 32 | 70 |
| F-VB-UST-9-14 | 240-57597-9 | 71 | 68 | 68 | 73 | 45 | 87 |
| F-VB-UST-10-14 | 240-57597-10 | 59 | 61 | 60 | 64 | 40 | 78 |
| F-VB-UST-11-14 | 240-57597-11 | 78 | 80 | 80 | 73 | 40 | 88 |
| F-VB-UST-12-14 | 240-57597-12 | 67 | 64 | 67 | 63 | 32 | 72 |
| F-VB-UST-13-14 | 240-57597-13 | 65 | 57 | 65 | 59 | 33 | 77 |
| F-VB-UST-14-14 | 240-57597-14 | 70 | 62 | 73 | 68 | 28 | 79 |
| F-VB-UST-DUP | 240-57597-15 | 71 | 60 | 74 | 68 | 24 | 81 |
| | MB 240-205786/23-A | 66 | 66 | 65 | 68 | 48 | 80 |
| | MB 240-205866/8-A | 81 | 88 | 78 | 76 | 49 | 87 |
| | LCS 240-205786/24-A | 79 | 79 | 79 | 81 | 44 | 93 |
| | LCS 240-205866/9-A | 72 | 75 | 76 | 64 | 33 | 76 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 69 | 62 | 78 | 71 | 30 | 80 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 61 | 53 | 65 | 60 | 36 | 77 |

| | QC LIMITS |
|---|---|
| 2FP = 2-Fluorophenol (Surr) | 24-110 |
| PHL = Phenol-d5 (Surr) | 26-110 |
| NBZ = Nitrobenzene-d5 (Surr) | 20-110 |
| FBP = 2-Fluorobiphenyl (Surr) | 24-110 |
| TBP = 2,4,6-Tribromophenol (Surr) | 10-110 |
| TPH = Terphenyl-d14 (Surr) | 36-110 |

# Column to be used to flag recovery values

FORM II 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111005.D

Lab ID: LCS 240-205786/24-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 525 | 79 | 35-110 | |
| bis (2-chloroisopropyl) ether | 667 | 611 | 92 | 29-110 | |
| 2,4,5-Trichlorophenol | 667 | 428 | 64 | 25-110 | |
| 2,4,6-Trichlorophenol | 667 | 319 | 48 | 12-110 | |
| 2,4-Dichlorophenol | 667 | 532 | 80 | 39-110 | |
| 2,4-Dimethylphenol | 667 | 462 | 69 | 29-110 | |
| 2,4-Dinitrophenol | 1330 | 481 | 36 | 10-110 | |
| 2,4-Dinitrotoluene | 667 | 605 | 91 | 48-110 | |
| 2,6-Dinitrotoluene | 667 | 563 | 84 | 45-110 | |
| 2-Chloronaphthalene | 667 | 523 | 78 | 32-110 | |
| 2-Chlorophenol | 667 | 528 | 79 | 37-110 | |
| 2-Methylnaphthalene | 667 | 550 | 83 | 36-110 | |
| 2-Methylphenol | 667 | 576 | 86 | 41-110 | |
| 2-Nitroaniline | 667 | 596 | 89 | 45-110 | |
| 2-Nitrophenol | 667 | 584 | 88 | 34-110 | |
| 3,3'-Dichlorobenzidine | 1330 | 900 | 68 | 28-110 | |
| 3-Nitroaniline | 667 | 497 | 75 | 44-110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 716 | 54 | 10-110 | |
| 4-Bromophenyl phenyl ether | 667 | 571 | 86 | 39-110 | |
| 4-Chloro-3-methylphenol | 667 | 571 | 86 | 48-110 | |
| 4-Chloroaniline | 667 | 395 | 59 | 30-110 | |
| 4-Chlorophenyl phenyl ether | 667 | 549 | 82 | 40-110 | |
| 4-Nitroaniline | 667 | 588 | 88 | 48-110 | |
| 4-Nitrophenol | 1330 | 1060 | 80 | 28-110 | |
| Acenaphthene | 667 | 533 | 80 | 38-110 | |
| Acenaphthylene | 667 | 532 | 80 | 40-110 | |
| Acetophenone | 667 | 551 | 83 | 40-110 | |
| Anthracene | 667 | 545 | 82 | 48-110 | |
| Atrazine | 1330 | 1240 | 93 | 54-120 | |
| Benzaldehyde | 1330 | 1150 | 86 | 32-110 | |
| Benzo[a]anthracene | 667 | 553 | 83 | 50-110 | |
| Benzo[a]pyrene | 667 | 592 | 89 | 44-110 | |
| Benzo[b]fluoranthene | 667 | 592 | 89 | 43-110 | |
| Benzo[g,h,i]perylene | 667 | 573 | 86 | 51-110 | |
| Benzo[k]fluoranthene | 667 | 571 | 86 | 38-105 | |
| Bis(2-chloroethoxy)methane | 667 | 537 | 80 | 32-110 | |
| Bis(2-chloroethyl)ether | 667 | 533 | 80 | 34-110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 662 | 99 | 50-110 | |
| Butyl benzyl phthalate | 667 | 604 | 91 | 51-110 | |
| Caprolactam | 1330 | 1110 | 83 | 44-114 | |
| Carbazole | 667 | 564 | 85 | 50-110 | |
| Chrysene | 667 | 550 | 83 | 50-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111005.D

Lab ID: LCS 240-205786/24-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 667 | 620 | 93 | 51-110 | |
| Dibenzofuran | 667 | 548 | 82 | 43-110 | |
| Diethyl phthalate | 667 | 618 | 93 | 52-110 | |
| Dimethyl phthalate | 667 | 572 | 86 | 50-110 | |
| Di-n-butyl phthalate | 667 | 583 | 88 | 51-110 | |
| Di-n-octyl phthalate | 667 | 628 | 94 | 48-110 | |
| Fluoranthene | 667 | 539 | 81 | 51-110 | |
| Fluorene | 667 | 550 | 83 | 46-110 | |
| Hexachlorobenzene | 667 | 553 | 83 | 43-110 | |
| Hexachlorobutadiene | 667 | 559 | 84 | 29-110 | |
| Hexachlorocyclopentadiene | 667 | 240 J | 36 | 12-110 | |
| Hexachloroethane | 667 | 539 | 81 | 30-110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 606 | 91 | 50-110 | |
| Isophorone | 667 | 547 | 82 | 36-110 | |
| N-Nitrosodi-n-propylamine | 667 | 563 | 84 | 38-110 | |
| N-Nitrosodiphenylamine | 1330 | 1180 | 89 | 46-110 | |
| Naphthalene | 667 | 537 | 81 | 36-110 | |
| Nitrobenzene | 667 | 550 | 83 | 32-110 | |
| Pentachlorophenol | 1330 | 584 | 44 | 10-110 | |
| Phenanthrene | 667 | 547 | 82 | 49-110 | |
| Phenol | 667 | 542 | 81 | 38-110 | |
| Pyrene | 667 | 574 | 86 | 49-110 | |
| 3 & 4 Methylphenol | 667 | 526 | 79 | 40-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111007.D

Lab ID: LCS 240-205866/9-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 396 | 59 | 35-110 | |
| bis (2-chloroisopropyl) ether | 667 | 514 | 77 | 29-110 | |
| 2,4,5-Trichlorophenol | 667 | 325 | 49 | 25-110 | |
| 2,4,6-Trichlorophenol | 667 | 243 | 37 | 12-110 | |
| 2,4-Dichlorophenol | 667 | 446 | 67 | 39-110 | |
| 2,4-Dimethylphenol | 667 | 444 | 67 | 29-110 | |
| 2,4-Dinitrophenol | 1330 | 464 | 35 | 10-110 | |
| 2,4-Dinitrotoluene | 667 | 473 | 71 | 48-110 | |
| 2,6-Dinitrotoluene | 667 | 462 | 69 | 45-110 | |
| 2-Chloronaphthalene | 667 | 425 | 64 | 32-110 | |
| 2-Chlorophenol | 667 | 439 | 66 | 37-110 | |
| 2-Methylnaphthalene | 667 | 428 | 64 | 36-110 | |
| 2-Methylphenol | 667 | 438 | 66 | 41-110 | |
| 2-Nitroaniline | 667 | 548 | 82 | 45-110 | |
| 2-Nitrophenol | 667 | 459 | 69 | 34-110 | |
| 3,3'-Dichlorobenzidine | 1330 | 946 | 71 | 28-110 | |
| 3-Nitroaniline | 667 | 446 | 67 | 44-110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 724 | 54 | 10-110 | |
| 4-Bromophenyl phenyl ether | 667 | 468 | 70 | 39-110 | |
| 4-Chloro-3-methylphenol | 667 | 444 | 67 | 48-110 | |
| 4-Chloroaniline | 667 | 366 | 55 | 30-110 | |
| 4-Chlorophenyl phenyl ether | 667 | 412 | 62 | 40-110 | |
| 4-Nitroaniline | 667 | 452 | 68 | 48-110 | |
| 4-Nitrophenol | 1330 | 804 | 60 | 28-110 | |
| Acenaphthene | 667 | 386 | 58 | 38-110 | |
| Acenaphthylene | 667 | 400 | 60 | 40-110 | |
| Acetophenone | 667 | 424 | 64 | 40-110 | |
| Anthracene | 667 | 462 | 69 | 48-110 | |
| Atrazine | 1330 | 1140 | 86 | 54-120 | |
| Benzaldehyde | 1330 | 1090 | 82 | 32-110 | |
| Benzo[a]anthracene | 667 | 451 | 68 | 50-110 | |
| Benzo[a]pyrene | 667 | 491 | 74 | 44-110 | |
| Benzo[b]fluoranthene | 667 | 559 | 84 | 43-110 | |
| Benzo[g,h,i]perylene | 667 | 467 | 70 | 51-110 | |
| Benzo[k]fluoranthene | 667 | 429 | 64 | 38-105 | |
| Bis(2-chloroethoxy)methane | 667 | 489 | 73 | 32-110 | |
| Bis(2-chloroethyl)ether | 667 | 503 | 75 | 34-110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 544 | 82 | 50-110 | |
| Butyl benzyl phthalate | 667 | 512 | 77 | 51-110 | |
| Caprolactam | 1330 | 1000 | 75 | 44-114 | |
| Carbazole | 667 | 536 | 80 | 50-110 | |
| Chrysene | 667 | 442 | 66 | 50-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111007.D

Lab ID: LCS 240-205866/9-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 667 | 513 | 77 | 51-110 | |
| Dibenzofuran | 667 | 406 | 61 | 43-110 | |
| Diethyl phthalate | 667 | 418 | 63 | 52-110 | |
| Dimethyl phthalate | 667 | 455 | 68 | 50-110 | |
| Di-n-butyl phthalate | 667 | 538 | 81 | 51-110 | |
| Di-n-octyl phthalate | 667 | 448 | 67 | 48-110 | |
| Fluoranthene | 667 | 500 | 75 | 51-110 | |
| Fluorene | 667 | 426 | 64 | 46-110 | |
| Hexachlorobenzene | 667 | 496 | 74 | 43-110 | |
| Hexachlorobutadiene | 667 | 392 | 59 | 29-110 | |
| Hexachlorocyclopentadiene | 667 | 244 J | 37 | 12-110 | |
| Hexachloroethane | 667 | 423 | 63 | 30-110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 505 | 76 | 50-110 | |
| Isophorone | 667 | 513 | 77 | 36-110 | |
| N-Nitrosodi-n-propylamine | 667 | 592 | 89 | 38-110 | |
| N-Nitrosodiphenylamine | 1330 | 987 | 74 | 46-110 | |
| Naphthalene | 667 | 387 | 58 | 36-110 | |
| Nitrobenzene | 667 | 506 | 76 | 32-110 | |
| Pentachlorophenol | 1330 | 573 | 43 | 10-110 | |
| Phenanthrene | 667 | 439 | 66 | 49-110 | |
| Phenol | 667 | 537 | 81 | 38-110 | |
| Pyrene | 667 | 481 | 72 | 49-110 | |
| 3 & 4 Methylphenol | 667 | 432 | 65 | 40-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 51111024.D

Lab ID: 240-57597-15 MS  Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 778 | 4.1 U | 511 | 66 | 32-110 | |
| bis (2-chloroisopropyl) ether | 778 | 11 U | 646 | 83 | 11-110 | |
| 2,4,5-Trichlorophenol | 778 | 29 U | 290 | 37 | 10-117 | |
| 2,4,6-Trichlorophenol | 778 | 10 U | 318 | 41 | 10-110 | |
| 2,4-Dichlorophenol | 778 | 23 U | 444 | 57 | 10-110 | |
| 2,4-Dimethylphenol | 778 | 23 U | 275 | 35 | 10-110 | |
| 2,4-Dinitrophenol | 1560 | 24 U | 940 | 60 | 10-110 | |
| 2,4-Dinitrotoluene | 778 | 20 U | 616 | 79 | 32-110 | |
| 2,6-Dinitrotoluene | 778 | 24 U | 597 | 77 | 35-110 | |
| 2-Chloronaphthalene | 778 | 0.52 U | 546 | 70 | 28-110 | |
| 2-Chlorophenol | 778 | 9.6 U | 504 | 65 | 10-110 | |
| 2-Methylnaphthalene | 778 | 0.58 U | 528 | 68 | 10-133 | |
| 2-Methylphenol | 778 | 13 U | 391 | 50 | 24-110 | |
| 2-Nitroaniline | 778 | 11 U | 709 | 91 | 39-110 | |
| 2-Nitrophenol | 778 | 9.7 U | 593 | 76 | 10-110 | |
| 3,3'-Dichlorobenzidine | 1560 | 21 U | 1180 | 76 | 10-110 | |
| 3-Nitroaniline | 778 | 19 U | 533 | 68 | 10-110 | |
| 4,6-Dinitro-2-methylphenol | 1560 | 11 U | 1080 | 69 | 10-110 | |
| 4-Bromophenyl phenyl ether | 778 | 15 U | 592 | 76 | 33-110 | |
| 4-Chloro-3-methylphenol | 778 | 24 U | 231 | 30 | 25-110 | |
| 4-Chloroaniline | 778 | 20 U | 391 | 50 | 10-110 | |
| 4-Chlorophenyl phenyl ether | 778 | 15 U | 537 | 69 | 32-110 | |
| 4-Nitroaniline | 778 | 30 U | 570 | 73 | 10-110 | |
| 4-Nitrophenol | 1560 | 20 U | 383 J | 25 | 10-113 | |
| Acenaphthene | 778 | 0.89 U | 501 | 64 | 22-110 | |
| Acenaphthylene | 778 | 0.41 U | 518 | 67 | 24-110 | |
| Acetophenone | 778 | 11 U | 542 | 70 | 31-110 | |
| Anthracene | 778 | 0.91 U | 585 | 75 | 20-110 | |
| Atrazine | 1560 | 11 U | 1450 | 93 | 45-118 | |
| Benzaldehyde | 1560 | 14 U | 1350 | 87 | 23-110 | |
| Benzo[a]anthracene | 778 | 0.73 U | 556 | 71 | 10-122 | |
| Benzo[a]pyrene | 778 | 0.75 U | 623 | 80 | 10-110 | |
| Benzo[b]fluoranthene | 778 | 0.69 U | 673 | 87 | 12-118 | |
| Benzo[g,h,i]perylene | 778 | 0.41 U | 598 | 77 | 10-117 | |
| Benzo[k]fluoranthene | 778 | 0.79 U | 562 | 72 | 10-121 | |
| Bis(2-chloroethoxy)methane | 778 | 26 U | 611 | 78 | 26-110 | |
| Bis(2-chloroethyl)ether | 778 | 2.3 U | 607 | 78 | 21-110 | |
| Bis(2-ethylhexyl) phthalate | 778 | 22 J | 679 | 87 | 40-110 | |
| Butyl benzyl phthalate | 778 | 12 U | 649 | 83 | 44-110 | |
| Caprolactam | 1560 | 43 U | 1260 | 81 | 10-134 | |
| Carbazole | 778 | 31 U | 666 | 86 | 34-110 | |
| Chrysene | 778 | 1.3 U | 544 | 70 | 10-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111024.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 778 | 0.77 U | 637 | 82 | 14-113 | |
| Dibenzofuran | 778 | 0.77 U | 536 | 69 | 29-110 | |
| Diethyl phthalate | 778 | 19 U | 546 | 70 | 42-110 | |
| Dimethyl phthalate | 778 | 20 U | 595 | 76 | 41-110 | |
| Di-n-butyl phthalate | 778 | 17 U | 681 | 88 | 43-110 | |
| Di-n-octyl phthalate | 778 | 9.2 U | 559 | 72 | 24-119 | |
| Fluoranthene | 778 | 0.64 U | 621 | 80 | 10-110 | |
| Fluorene | 778 | 0.62 U | 548 | 70 | 23-110 | |
| Hexachlorobenzene | 778 | 2.4 U | 623 | 80 | 34-110 | |
| Hexachlorobutadiene | 778 | 6.5 U | 472 | 61 | 25-110 | |
| Hexachlorocyclopentadiene | 778 | 9.4 U | 219 J | 28 | 10-110 | |
| Hexachloroethane | 778 | 10 U | 526 | 68 | 12-110 | |
| Indeno[1,2,3-cd]pyrene | 778 | 0.41 U | 622 | 80 | 10-114 | |
| Isophorone | 778 | 15 U | 651 | 84 | 29-110 | |
| N-Nitrosodi-n-propylamine | 778 | 7.3 U | 757 | 97 | 26-110 | |
| N-Nitrosodiphenylamine | 1560 | 24 U | 1270 | 81 | 22-110 | |
| Naphthalene | 778 | 0.96 U | 473 | 61 | 10-111 | |
| Nitrobenzene | 778 | 2.6 U | 633 | 81 | 23-110 | |
| Pentachlorophenol | 1560 | 11 U | 423 | 27 | 10-110 | |
| Phenanthrene | 778 | 0.85 U | 557 | 72 | 10-166 | |
| Phenol | 778 | 8.5 U | 395 | 51 | 17-110 | |
| Pyrene | 778 | 0.51 U | 603 | 77 | 10-147 | |
| 3 & 4 Methylphenol | 778 | 23 U | 368 J | 47 | 25-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111025.D

Lab ID: 240-57597-15 MSD Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 781 | 419 | 54 | 20 | 32 | 32-110 | |
| bis (2-chloroisopropyl) ether | 781 | 536 | 69 | 19 | 42 | 11-110 | |
| 2,4,5-Trichlorophenol | 781 | 208 | 27 | 33 | 99 | 10-117 | |
| 2,4,6-Trichlorophenol | 781 | 307 | 39 | 4 | 38 | 10-110 | |
| 2,4-Dichlorophenol | 781 | 368 | 47 | 19 | 34 | 10-110 | |
| 2,4-Dimethylphenol | 781 | 286 | 37 | 4 | 31 | 10-110 | |
| 2,4-Dinitrophenol | 1560 | 819 | 52 | 14 | 99 | 10-110 | |
| 2,4-Dinitrotoluene | 781 | 589 | 75 | 4 | 30 | 32-110 | |
| 2,6-Dinitrotoluene | 781 | 546 | 70 | 9 | 30 | 35-110 | |
| 2-Chloronaphthalene | 781 | 455 | 58 | 18 | 30 | 28-110 | |
| 2-Chlorophenol | 781 | 422 | 54 | 18 | 47 | 10-110 | |
| 2-Methylnaphthalene | 781 | 437 | 56 | 19 | 42 | 10-133 | |
| 2-Methylphenol | 781 | 416 | 53 | 6 | 51 | 24-110 | |
| 2-Nitroaniline | 781 | 646 | 83 | 9 | 31 | 39-110 | |
| 2-Nitrophenol | 781 | 492 | 63 | 19 | 49 | 10-110 | |
| 3,3'-Dichlorobenzidine | 1560 | 1210 | 77 | 2 | 56 | 10-110 | |
| 3-Nitroaniline | 781 | 551 | 71 | 3 | 30 | 10-110 | |
| 4,6-Dinitro-2-methylphenol | 1560 | 996 | 64 | 8 | 55 | 10-110 | |
| 4-Bromophenyl phenyl ether | 781 | 532 | 68 | 11 | 30 | 33-110 | |
| 4-Chloro-3-methylphenol | 781 | 272 | 35 | 16 | 54 | 25-110 | |
| 4-Chloroaniline | 781 | 392 | 50 | 0 | 36 | 10-110 | |
| 4-Chlorophenyl phenyl ether | 781 | 476 | 61 | 12 | 30 | 32-110 | |
| 4-Nitroaniline | 781 | 541 | 69 | 5 | 48 | 10-110 | |
| 4-Nitrophenol | 1560 | 446 | 29 | 15 | 49 | 10-113 | |
| Acenaphthene | 781 | 432 | 55 | 15 | 99 | 22-110 | |
| Acenaphthylene | 781 | 443 | 57 | 16 | 99 | 24-110 | |
| Acetophenone | 781 | 442 | 57 | 20 | 43 | 31-110 | |
| Anthracene | 781 | 523 | 67 | 11 | 99 | 20-110 | |
| Atrazine | 1560 | 1300 | 83 | 11 | 30 | 45-118 | |
| Benzaldehyde | 1560 | 1100 | 71 | 20 | 42 | 23-110 | |
| Benzo[a]anthracene | 781 | 536 | 69 | 4 | 99 | 10-122 | |
| Benzo[a]pyrene | 781 | 563 | 72 | 10 | 99 | 10-110 | |
| Benzo[b]fluoranthene | 781 | 638 | 82 | 5 | 99 | 12-118 | |
| Benzo[g,h,i]perylene | 781 | 545 | 70 | 9 | 99 | 10-117 | |
| Benzo[k]fluoranthene | 781 | 486 | 62 | 14 | 99 | 10-121 | |
| Bis(2-chloroethoxy)methane | 781 | 497 | 64 | 21 | 37 | 26-110 | |
| Bis(2-chloroethyl)ether | 781 | 528 | 68 | 14 | 55 | 21-110 | |
| Bis(2-ethylhexyl) phthalate | 781 | 661 | 85 | 3 | 30 | 40-110 | |
| Butyl benzyl phthalate | 781 | 614 | 79 | 5 | 30 | 44-110 | |
| Caprolactam | 1560 | 1170 | 75 | 8 | 32 | 10-134 | |
| Carbazole | 781 | 604 | 77 | 10 | 30 | 34-110 | |
| Chrysene | 781 | 521 | 67 | 4 | 99 | 10-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 51111025.D

Lab ID: 240-57597-15 MSD  Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 781 | 596 | 76 | 7 | 99 | 14-113 | |
| Dibenzofuran | 781 | 463 | 59 | 15 | 30 | 29-110 | |
| Diethyl phthalate | 781 | 507 | 65 | 7 | 30 | 42-110 | |
| Dimethyl phthalate | 781 | 543 | 69 | 9 | 30 | 41-110 | |
| Di-n-butyl phthalate | 781 | 631 | 81 | 8 | 30 | 43-110 | |
| Di-n-octyl phthalate | 781 | 517 | 66 | 8 | 30 | 24-119 | |
| Fluoranthene | 781 | 577 | 74 | 7 | 99 | 10-110 | |
| Fluorene | 781 | 486 | 62 | 12 | 99 | 23-110 | |
| Hexachlorobenzene | 781 | 557 | 71 | 11 | 30 | 34-110 | |
| Hexachlorobutadiene | 781 | 395 | 51 | 18 | 34 | 25-110 | |
| Hexachlorocyclopentadiene | 781 | 186 J | 24 | 16 | 79 | 10-110 | |
| Hexachloroethane | 781 | 411 | 53 | 25 | 50 | 12-110 | |
| Indeno[1,2,3-cd]pyrene | 781 | 578 | 74 | 7 | 99 | 10-114 | |
| Isophorone | 781 | 540 | 69 | 19 | 38 | 29-110 | |
| N-Nitrosodi-n-propylamine | 781 | 615 | 79 | 21 | 42 | 26-110 | |
| N-Nitrosodiphenylamine | 1560 | 1130 | 72 | 11 | 30 | 22-110 | |
| Naphthalene | 781 | 397 | 51 | 17 | 99 | 10-111 | |
| Nitrobenzene | 781 | 515 | 66 | 21 | 41 | 23-110 | |
| Pentachlorophenol | 1560 | 617 | 39 | 37 | 50 | 10-110 | |
| Phenanthrene | 781 | 510 | 65 | 9 | 99 | 10-166 | |
| Phenol | 781 | 418 | 53 | 6 | 53 | 17-110 | |
| Pyrene | 781 | 575 | 74 | 5 | 99 | 10-147 | |
| 3 & 4 Methylphenol | 781 | 392 J | 50 | 6 | 50 | 25-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM IV
GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab File ID: 51111004.D  Lab Sample ID: MB 240-205786/23-A

Matrix: Solid  Date Extracted: 11/09/2015 06:12

Instrument ID: A4HP10  Date Analyzed: 11/11/2015 10:27

Level:(Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-205786/24-A | 51111005.D | 11/11/2015 10:51 |
| F-VB-UST-1-14 | 240-57597-1 | 51111006.D | 11/11/2015 11:16 |
| F-VB-UST-2-14 | 240-57597-2 | 51111007.D | 11/11/2015 11:40 |
| F-VB-UST-3-14 | 240-57597-3 | 51111008.D | 11/11/2015 12:05 |
| F-VB-UST-4-14 | 240-57597-4 | 51111009.D | 11/11/2015 12:30 |
| F-VB-UST-5-14 | 240-57597-5 | 51111010.D | 11/11/2015 12:54 |
| F-VB-UST-6-14 | 240-57597-6 | 51111011.D | 11/11/2015 13:19 |
| F-VB-UST-7-14 | 240-57597-7 | 51111012.D | 11/11/2015 13:44 |
| F-VB-UST-8-14 | 240-57597-8 | 51111013.D | 11/11/2015 14:08 |
| F-VB-UST-9-14 | 240-57597-9 | 51111014.D | 11/11/2015 14:33 |
| F-VB-UST-10-14 | 240-57597-10 | 51111015.D | 11/11/2015 14:57 |

FORM IV
GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab File ID: 51111006.D  Lab Sample ID: MB 240-205866/8-A

Matrix: Solid  Date Extracted: 11/09/2015 12:23

Instrument ID: A4HP9  Date Analyzed: 11/11/2015 09:54

Level:(Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-205866/9-A | 51111007.D | 11/11/2015 10:19 |
| F-VB-UST-11-14 | 240-57597-11 | 51111019.D | 11/11/2015 14:50 |
| F-VB-UST-12-14 | 240-57597-12 | 51111020.D | 11/11/2015 15:14 |
| F-VB-UST-13-14 | 240-57597-13 | 51111021.D | 11/11/2015 15:39 |
| F-VB-UST-14-14 | 240-57597-14 | 51111022.D | 11/11/2015 16:03 |
| F-VB-UST-DUP | 240-57597-15 | 51111023.D | 11/11/2015 16:28 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 51111024.D | 11/11/2015 16:52 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 51111025.D | 11/11/2015 17:17 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 51002001.D DFTPP Injection Date: 10/02/2015

Instrument ID: A4HP10 DFTPP Injection Time: 10:22

Analysis Batch No.: 200091

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 41.5 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0)1 |
| 69 | Mass 69 relative abundance | 35.2 | |
| 70 | Less than 2.0% of mass 69 | 0.2 | (0.7)1 |
| 127 | 25.0 - 75.0% of mass 198 | 49.0 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.8 | |
| 275 | 10.0- 30.0% of mass 198 | 25.4 | |
| 365 | Greater than 0.75% of mass 198 | 3.2 | |
| 441 | Present, but less than mass 443 | 16.6 | |
| 442 | 40.0 - 110.0% of mass 198 | 102.9 | |
| 443 | 15.0 - 24.0% of mass 442 | 20.5 | (20.0)2 |

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD5 240-200091/2 | 51002002.D | 10/02/2015 | 10:48 |
| | STD4 240-200091/3 | 51002003.D | 10/02/2015 | 11:12 |
| | STD3 240-200091/4 | 51002004.D | 10/02/2015 | 11:37 |
| | STD2 240-200091/5 | 51002005.D | 10/02/2015 | 12:01 |
| | STD1 240-200091/6 | 51002006.D | 10/02/2015 | 12:25 |
| | STD9 240-200091/7 | 51002007.D | 10/02/2015 | 12:50 |
| | STD8 240-200091/8 | 51002008.D | 10/02/2015 | 13:15 |
| | STD7 240-200091/9 | 51002009.D | 10/02/2015 | 13:39 |
| | STD6 240-200091/10 | 51002010.D | 10/02/2015 | 14:04 |
| | ICV 240-200091/11 | 51002011.D | 10/02/2015 | 14:28 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 51111001.D DFTPP Injection Date: 11/11/2015

Instrument ID: A4HP10 DFTPP Injection Time: 09:12

Analysis Batch No.: 206256

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 40.9 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0)1 |
| 69 | Mass 69 relative abundance | 34.5 | |
| 70 | Less than 2.0% of mass 69 | 0.2 | (0.5)1 |
| 127 | 25.0 - 75.0% of mass 198 | 50.8 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.5 | |
| 275 | 10.0- 30.0% of mass 198 | 25.4 | |
| 365 | Greater than 0.75% of mass 198 | 3.2 | |
| 441 | Present, but less than mass 443 | 8.7 | |
| 442 | 40.0 - 110.0% of mass 198 | 95.2 | |
| 443 | 15.0 - 24.0% of mass 442 | 18.5 | (19.4)2 |

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCV 240-206256/2 | 51111002.D | 11/11/2015 | 09:34 |
| | MB 240-205786/23-A | 51111004.D | 11/11/2015 | 10:27 |
| | LCS 240-205786/24-A | 51111005.D | 11/11/2015 | 10:51 |
| F-VB-UST-1-14 | 240-57597-1 | 51111006.D | 11/11/2015 | 11:16 |
| F-VB-UST-2-14 | 240-57597-2 | 51111007.D | 11/11/2015 | 11:40 |
| F-VB-UST-3-14 | 240-57597-3 | 51111008.D | 11/11/2015 | 12:05 |
| F-VB-UST-4-14 | 240-57597-4 | 51111009.D | 11/11/2015 | 12:30 |
| F-VB-UST-5-14 | 240-57597-5 | 51111010.D | 11/11/2015 | 12:54 |
| F-VB-UST-6-14 | 240-57597-6 | 51111011.D | 11/11/2015 | 13:19 |
| F-VB-UST-7-14 | 240-57597-7 | 51111012.D | 11/11/2015 | 13:44 |
| F-VB-UST-8-14 | 240-57597-8 | 51111013.D | 11/11/2015 | 14:08 |
| F-VB-UST-9-14 | 240-57597-9 | 51111014.D | 11/11/2015 | 14:33 |
| F-VB-UST-10-14 | 240-57597-10 | 51111015.D | 11/11/2015 | 14:57 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 51030001.D DFTPP Injection Date: 10/30/2015

Instrument ID: A4HP9 DFTPP Injection Time: 12:57

Analysis Batch No.: 204569

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 46.3 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0)1 |
| 69 | Mass 69 relative abundance | 51.4 | |
| 70 | Less than 2.0% of mass 69 | 0.2 | (0.4)1 |
| 127 | 25.0 - 75.0% of mass 198 | 59.4 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.3 | |
| 275 | 10.0- 30.0% of mass 198 | 21.2 | |
| 365 | Greater than 0.75% of mass 198 | 3.2 | |
| 441 | Present, but less than mass 443 | 0.3 | |
| 442 | 40.0 - 110.0% of mass 198 | 85.4 | |
| 443 | 15.0 - 24.0% of mass 442 | 16.8 | (19.7)2 |

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD5 240-204569/2 | 51030002.D | 10/30/2015 | 13:19 |
| | STD4 240-204569/3 | 51030003.D | 10/30/2015 | 13:44 |
| | STD3 240-204569/4 | 51030004.D | 10/30/2015 | 14:10 |
| | STD2 240-204569/5 | 51030005.D | 10/30/2015 | 14:35 |
| | STD1 240-204569/6 | 51030006.D | 10/30/2015 | 15:01 |
| | STD9 240-204569/7 | 51030007.D | 10/30/2015 | 15:27 |
| | STD8 240-204569/8 | 51030008.D | 10/30/2015 | 15:52 |
| | STD7 240-204569/9 | 51030009.D | 10/30/2015 | 16:17 |
| | STD6 240-204569/10 | 51030010.D | 10/30/2015 | 16:42 |
| | ICV 240-204569/11 | 51030011.D | 10/30/2015 | 17:08 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 51111001.D DFTPP Injection Date: 11/11/2015

Instrument ID: A4HP9 DFTPP Injection Time: 07:27

Analysis Batch No.: 206233

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 50.3 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0)1 |
| 69 | Mass 69 relative abundance | 58.1 | |
| 70 | Less than 2.0% of mass 69 | 0.3 | (0.5)1 |
| 127 | 25.0 - 75.0% of mass 198 | 54.3 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.6 | |
| 275 | 10.0- 30.0% of mass 198 | 22.9 | |
| 365 | Greater than 0.75% of mass 198 | 2.7 | |
| 441 | Present, but less than mass 443 | 10.7 | |
| 442 | 40.0 - 110.0% of mass 198 | 61.4 | |
| 443 | 15.0 - 24.0% of mass 442 | 11.9 | (19.4)2 |

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCV 240-206233/2 | 51111002.D | 11/11/2015 | 07:49 |
| | MB 240-205866/8-A | 51111006.D | 11/11/2015 | 09:54 |
| | LCS 240-205866/9-A | 51111007.D | 11/11/2015 | 10:19 |
| F-VB-UST-11-14 | 240-57597-11 | 51111019.D | 11/11/2015 | 14:50 |
| F-VB-UST-12-14 | 240-57597-12 | 51111020.D | 11/11/2015 | 15:14 |
| F-VB-UST-13-14 | 240-57597-13 | 51111021.D | 11/11/2015 | 15:39 |
| F-VB-UST-14-14 | 240-57597-14 | 51111022.D | 11/11/2015 | 16:03 |
| F-VB-UST-DUP | 240-57597-15 | 51111023.D | 11/11/2015 | 16:28 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 51111024.D | 11/11/2015 | 16:52 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 51111025.D | 11/11/2015 | 17:17 |

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Sample No.: STD6 240-200091/10 Date Analyzed: 10/02/2015 14:04

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm)

Lab File ID (Standard): 51002010.D Heated Purge: (Y/N) N

Calibration ID: 30980

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 221463 | 6.23 | 906472 | 7.34 | 529260 | 8.85 |
| UPPER LIMIT | | 442926 | 6.73 | 1812944 | 7.84 | 1058520 | 9.35 |
| LOWER LIMIT | | 110732 | 5.73 | 453236 | 6.84 | 264630 | 8.35 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-200091/11 | | 242280 | 6.23 | 1000939 | 7.34 | 580042 | 8.85 |
| CCV 240-206256/2 CCVIS | | 183138 | 6.15 | 734641 | 7.27 | 425153 | 8.77 |

DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Sample No.: STD6 240-200091/10 Date Analyzed: 10/02/2015 14:04

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm)

Lab File ID (Standard): 51002010.D Heated Purge: (Y/N) N

Calibration ID: 30980

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 840219 | 10.12 | 1037091 | 12.82 | 955028 | 15.02 |
| UPPER LIMIT | | 1680438 | 10.62 | 2074182 | 13.32 | 1910056 | 15.52 |
| LOWER LIMIT | | 420110 | 9.62 | 518546 | 12.32 | 477514 | 14.52 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-200091/11 | | 929663 | 10.12 | 1154414 | 12.82 | 1074213 | 15.02 |
| CCV 240-206256/2 CCVIS | | 678359 | 10.04 | 794127 | 12.71 | 678609 | 14.89 |

PHN = Phenanthrene-d10
CRY = Chrysene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206256/2  Date Analyzed: 11/11/2015 09:34

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 51111002.D  Heated Purge: (Y/N) N

Calibration ID: 31049

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 183138 | 6.15 | 734641 | 7.27 | 425153 | 8.77 |
| UPPER LIMIT | | 366276 | 6.65 | 1469282 | 7.77 | 850306 | 9.27 |
| LOWER LIMIT | | 91569 | 5.65 | 367321 | 6.77 | 212577 | 8.27 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-205786/23-A | | 169364 | 6.16 | 695817 | 7.27 | 416939 | 8.77 |
| LCS 240-205786/24-A | | 167619 | 6.16 | 683191 | 7.27 | 403392 | 8.77 |
| 240-57597-1 | F-VB-UST-1-14 | 170794 | 6.16 | 710718 | 7.27 | 416040 | 8.77 |
| 240-57597-2 | F-VB-UST-2-14 | 172301 | 6.16 | 704300 | 7.27 | 408420 | 8.77 |
| 240-57597-3 | F-VB-UST-3-14 | 168152 | 6.16 | 687534 | 7.27 | 389818 | 8.77 |
| 240-57597-4 | F-VB-UST-4-14 | 159703 | 6.16 | 681534 | 7.27 | 397658 | 8.77 |
| 240-57597-5 | F-VB-UST-5-14 | 169966 | 6.16 | 709398 | 7.27 | 411661 | 8.77 |
| 240-57597-6 | F-VB-UST-6-14 | 173464 | 6.16 | 694821 | 7.27 | 402882 | 8.77 |
| 240-57597-7 | F-VB-UST-7-14 | 166312 | 6.16 | 698152 | 7.27 | 395626 | 8.77 |
| 240-57597-8 | F-VB-UST-8-14 | 170392 | 6.16 | 690624 | 7.27 | 399583 | 8.77 |
| 240-57597-9 | F-VB-UST-9-14 | 167764 | 6.16 | 674344 | 7.27 | 396455 | 8.77 |
| 240-57597-10 | F-VB-UST-10-14 | 165361 | 6.16 | 676786 | 7.27 | 404905 | 8.77 |

DCB = 1,4-Dichlorobenzene-d4
DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
NPT = Naphthalene-d8
Area Limit = 50%-200% of internal standard area
ANT = Acenaphthene-d10
RT Limit = ± 0.5 minutes of internal standard RT
ANT = Acenaphthene-d10

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206256/2  Date Analyzed: 11/11/2015 09:34

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 51111002.D  Heated Purge: (Y/N) N

Calibration ID: 31049

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 678359 | 10.04 | 794127 | 12.71 | 678609 | 14.89 |
| UPPER LIMIT | | 1356718 | 10.54 | 1588254 | 13.21 | 1357218 | 15.39 |
| LOWER LIMIT | | 339180 | 9.54 | 397064 | 12.21 | 339305 | 14.39 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-205786/23-A | | 633596 | 10.04 | 730132 | 12.70 | 644324 | 14.88 |
| LCS 240-205786/24-A | | 651601 | 10.04 | 738522 | 12.70 | 635222 | 14.89 |
| 240-57597-1 | F-VB-UST-1-14 | 661549 | 10.04 | 736014 | 12.70 | 508559 | 14.88 |
| 240-57597-2 | F-VB-UST-2-14 | 651468 | 10.04 | 719477 | 12.70 | 573126 | 14.88 |
| 240-57597-3 | F-VB-UST-3-14 | 616636 | 10.04 | 695392 | 12.70 | 523351 | 14.88 |
| 240-57597-4 | F-VB-UST-4-14 | 624736 | 10.04 | 703819 | 12.70 | 478985 | 14.88 |
| 240-57597-5 | F-VB-UST-5-14 | 639217 | 10.04 | 741747 | 12.70 | 497306 | 14.88 |
| 240-57597-6 | F-VB-UST-6-14 | 640684 | 10.04 | 728874 | 12.70 | 488005 | 14.88 |
| 240-57597-7 | F-VB-UST-7-14 | 639827 | 10.04 | 723063 | 12.70 | 509629 | 14.88 |
| 240-57597-8 | F-VB-UST-8-14 | 631926 | 10.04 | 719274 | 12.70 | 575492 | 14.88 |
| 240-57597-9 | F-VB-UST-9-14 | 624386 | 10.04 | 704835 | 12.70 | 495093 | 14.88 |
| 240-57597-10 | F-VB-UST-10-14 | 640666 | 10.04 | 718652 | 12.70 | 538508 | 14.88 |

PHN = Phenanthrene-d10
PHN = Phenanthrene-d10
CRY = Chrysene-d12
CRY = Chrysene-d12
PRY = Perylene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Sample No.: STD6 240-204569/10 Date Analyzed: 10/30/2015 16:42

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm)

Lab File ID (Standard): 51030010.D Heated Purge: (Y/N) N

Calibration ID: 31170

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 251143 | 6.42 | 1074814 | 7.54 | 555801 | 9.04 |
| UPPER LIMIT | | 502286 | 6.92 | 2149628 | 8.04 | 1111602 | 9.54 |
| LOWER LIMIT | | 125572 | 5.92 | 537407 | 7.04 | 277901 | 8.54 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-204569/11 | | 291290 | 6.42 | 1233894 | 7.54 | 643573 | 9.04 |
| CCV 240-206233/2 CCVIS | | 263455 | 6.30 | 1198567 | 7.42 | 612244 | 8.92 |

DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Sample No.: STD6 240-204569/10 Date Analyzed: 10/30/2015 16:42

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm)

Lab File ID (Standard): 51030010.D Heated Purge: (Y/N) N

Calibration ID: 31170

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 848550 | 10.31 | 938909 | 12.62 | 798084 | 14.43 |
| UPPER LIMIT | | 1697100 | 10.81 | 1877818 | 13.12 | 1596168 | 14.93 |
| LOWER LIMIT | | 424275 | 9.81 | 469455 | 12.12 | 399042 | 13.93 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-204569/11 | | 1041455 | 10.31 | 1160345 | 12.61 | 991819 | 14.43 |
| CCV 240-206233/2 CCVIS | | 908264 | 10.19 | 962378 | 12.46 | 831716 | 14.18 |

PHN = Phenanthrene-d10
CRY = Chrysene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206233/2  Date Analyzed: 11/11/2015 07:49

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)

Lab File ID (Standard): 51111002.D  Heated Purge: (Y/N) N

Calibration ID: 31170

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 263455 | 6.30 | 1198567 | 7.42 | 612244 | 8.92 |
| UPPER LIMIT | | 526910 | 6.80 | 2397134 | 7.92 | 1224488 | 9.42 |
| LOWER LIMIT | | 131728 | 5.80 | 599284 | 6.92 | 306122 | 8.42 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-205866/8-A | | 217654 | 6.30 | 987205 | 7.42 | 480730 | 8.92 |
| LCS 240-205866/9-A | | 227686 | 6.30 | 1006110 | 7.42 | 522813 | 8.92 |
| 240-57597-11 | F-VB-UST-11-14 | 214487 | 6.29 | 930769 | 7.42 | 469347 | 8.92 |
| 240-57597-12 | F-VB-UST-12-14 | 207603 | 6.30 | 930370 | 7.42 | 469544 | 8.92 |
| 240-57597-13 | F-VB-UST-13-14 | 229993 | 6.30 | 1005170 | 7.42 | 509477 | 8.92 |
| 240-57597-14 | F-VB-UST-14-14 | 236787 | 6.30 | 998849 | 7.42 | 507858 | 8.92 |
| 240-57597-15 | F-VB-UST-DUP | 220994 | 6.30 | 932862 | 7.42 | 478231 | 8.92 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 216080 | 6.30 | 960244 | 7.42 | 481718 | 8.92 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 234420 | 6.30 | 1033463 | 7.42 | 522594 | 8.92 |

DCB = 1,4-Dichlorobenzene-d4
DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
NPT = Naphthalene-d8
Area Limit = 50%-200% of internal standard area
ANT = Acenaphthene-d10
RT Limit = ± 0.5 minutes of internal standard RT
ANT = Acenaphthene-d10

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206233/2  Date Analyzed: 11/11/2015 07:49

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)

Lab File ID (Standard): 51111002.D  Heated Purge: (Y/N) N

Calibration ID: 31170

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 908264 | 10.19 | 962378 | 12.46 | 831716 | 14.18 |
| UPPER LIMIT | | 1816528 | 10.69 | 1924756 | 12.96 | 1663432 | 14.68 |
| LOWER LIMIT | | 454132 | 9.69 | 481189 | 11.96 | 415858 | 13.68 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-205866/8-A | | 733495 | 10.18 | 742087 | 12.46 | 654968 | 14.18 |
| LCS 240-205866/9-A | | 774857 | 10.19 | 814640 | 12.46 | 725274 | 14.18 |
| 240-57597-11 | F-VB-UST-11-14 | 720283 | 10.19 | 731765 | 12.46 | 528934 | 14.18 |
| 240-57597-12 | F-VB-UST-12-14 | 701609 | 10.19 | 717786 | 12.46 | 585455 | 14.18 |
| 240-57597-13 | F-VB-UST-13-14 | 754672 | 10.19 | 785246 | 12.46 | 684137 | 14.18 |
| 240-57597-14 | F-VB-UST-14-14 | 760631 | 10.19 | 787874 | 12.46 | 628812 | 14.18 |
| 240-57597-15 | F-VB-UST-DUP | 704693 | 10.18 | 734178 | 12.46 | 582028 | 14.17 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 739261 | 10.19 | 786491 | 12.46 | 699060 | 14.18 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 822314 | 10.19 | 841298 | 12.46 | 788802 | 14.18 |

PHN = Phenanthrene-d10
PHN = Phenanthrene-d10
CRY = Chrysene-d12
CRY = Chrysene-d12
PRY = Perylene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 51111006.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.99(g) | Date Analyzed: 11/11/2015 11:16

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 59 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.53 | U | 59 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 9.7 | U | 59 | 9.7 |
| 91-57-6 | 2-Methylnaphthalene | 0.59 | U | 7.9 | 0.59 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.9 | U | 59 | 9.9 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 59 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 59 | 15 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.90 | U | 7.9 | 0.90 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 7.9 | 0.42 |
| 98-86-2 | Acetophenone | 16 | J | 120 | 11 |
| 120-12-7 | Anthracene | 0.93 | U | 7.9 | 0.93 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.75 | U | 7.9 | 0.75 |
| 50-32-8 | Benzo[a]pyrene | 0.76 | U | 7.9 | 0.76 |
| 205-99-2 | Benzo[b]fluoranthene | 0.70 | U | 7.9 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.42 | U | 7.9 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 51111006.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.99(g) | Date Analyzed: 11/11/2015 11:16

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.81 | U | 7.9 | 0.81 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 31 | J | 83 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 83 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 59 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 7.9 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.78 | U | 7.9 | 0.78 |
| 132-64-9 | Dibenzofuran | 0.78 | U | 59 | 0.78 |
| 84-66-2 | Diethyl phthalate | 19 | U | 83 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 83 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 83 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.4 | U | 83 | 9.4 |
| 206-44-0 | Fluoranthene | 0.65 | U | 7.9 | 0.65 |
| 86-73-7 | Fluorene | 0.63 | U | 7.9 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.9 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.6 | U | 59 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.6 | U | 390 | 9.6 |
| 67-72-1 | Hexachloroethane | 11 | U | 59 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.42 | U | 7.9 | 0.42 |
| 78-59-1 | Isophorone | 15 | U | 59 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.5 | U | 59 | 7.5 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 59 | 25 |
| 91-20-3 | Naphthalene | 0.97 | U | 7.9 | 0.97 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.87 | U | 7.9 | 0.87 |
| 108-95-2 | Phenol | 8.7 | U | 59 | 8.7 |
| 129-00-0 | Pyrene | 0.52 | U | 7.9 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 470 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 51111006.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.99(g) | Date Analyzed: 11/11/2015 11:16

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 76 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 65 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 62 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 64 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 67 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 44 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 51111007.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:02

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.16(g) | Date Analyzed: 11/11/2015 11:40

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 58 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 29 | U | 180 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10 | U | 180 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 23 | U | 180 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 23 | U | 180 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 230 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.53 | U | 58 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 9.6 | U | 58 | 9.6 |
| 91-57-6 | 2-Methylnaphthalene | 0.58 | U | 7.8 | 0.58 |
| 95-48-7 | 2-Methylphenol | 13 | U | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.7 | U | 58 | 9.7 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 58 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 58 | 15 |
| 100-01-6 | 4-Nitroaniline | 30 | U | 230 | 30 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.89 | U | 7.8 | 0.89 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.8 | 0.41 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.91 | U | 7.8 | 0.91 |
| 1912-24-9 | Atrazine | 11 | U | 230 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.74 | U | 7.8 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 0.75 | U | 7.8 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 0.69 | U | 7.8 | 0.69 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.41 | U | 7.8 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 51111007.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:02

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.16(g) | Date Analyzed: 11/11/2015 11:40

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.80 | U | 7.8 | 0.80 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.3 | U | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 22 | U | 82 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 82 | 12 |
| 105-60-2 | Caprolactam | 43 | U | 390 | 43 |
| 86-74-8 | Carbazole | 32 | U | 58 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 7.8 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.77 | U | 7.8 | 0.77 |
| 132-64-9 | Dibenzofuran | 0.77 | U | 58 | 0.77 |
| 84-66-2 | Diethyl phthalate | 19 | U | 82 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 82 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 82 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.2 | U | 82 | 9.2 |
| 206-44-0 | Fluoranthene | 0.64 | U | 7.8 | 0.64 |
| 86-73-7 | Fluorene | 0.62 | U | 7.8 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.8 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.6 | U | 58 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.5 | U | 390 | 9.5 |
| 67-72-1 | Hexachloroethane | 11 | U | 58 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.41 | U | 7.8 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 58 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.4 | U | 58 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 58 | 25 |
| 91-20-3 | Naphthalene | 0.96 | U | 7.8 | 0.96 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.85 | U | 7.8 | 0.85 |
| 108-95-2 | Phenol | 8.5 | U | 58 | 8.5 |
| 129-00-0 | Pyrene | 0.51 | U | 7.8 | 0.51 |
| 15831-10-4 | 3 & 4 Methylphenol | 23 | U | 470 | 23 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 51111007.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:02

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.16(g) | Date Analyzed: 11/11/2015 11:40

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 78 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 61 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 58 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 61 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 63 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 42 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 51111008.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:04

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.03(g) | Date Analyzed: 11/11/2015 12:05

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 17.8 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.3 | U | 61 | 4.3 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 12 | U | 120 | 12 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 26 | U | 400 | 26 |
| 121-14-2 | 2,4-Dinitrotoluene | 21 | U | 240 | 21 |
| 606-20-2 | 2,6-Dinitrotoluene | 26 | U | 240 | 26 |
| 91-58-7 | 2-Chloronaphthalene | 0.55 | U | 61 | 0.55 |
| 95-57-8 | 2-Chlorophenol | 10 | U | 61 | 10 |
| 91-57-6 | 2-Methylnaphthalene | 0.61 | U | 8.1 | 0.61 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 10 | U | 61 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 61 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 26 | U | 180 | 26 |
| 106-47-8 | 4-Chloroaniline | 21 | U | 180 | 21 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 61 | 16 |
| 100-01-6 | 4-Nitroaniline | 32 | U | 240 | 32 |
| 100-02-7 | 4-Nitrophenol | 21 | U | 400 | 21 |
| 83-32-9 | Acenaphthene | 0.92 | U | 8.1 | 0.92 |
| 208-96-8 | Acenaphthylene | 0.43 | U | 8.1 | 0.43 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.95 | U | 8.1 | 0.95 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 15 | U | 120 | 15 |
| 56-55-3 | Benzo[a]anthracene | 0.77 | U | 8.1 | 0.77 |
| 50-32-8 | Benzo[a]pyrene | 0.78 | U | 8.1 | 0.78 |
| 205-99-2 | Benzo[b]fluoranthene | 0.72 | U | 8.1 | 0.72 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.43 | U | 8.1 | 0.43 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 51111008.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:04

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.03(g) | Date Analyzed: 11/11/2015 12:05

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 17.8 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.83 | U | 8.1 | 0.83 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 27 | U | 120 | 27 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 23 | U | 85 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 85 | 12 |
| 105-60-2 | Caprolactam | 45 | U | 400 | 45 |
| 86-74-8 | Carbazole | 33 | U | 61 | 33 |
| 218-01-9 | Chrysene | 1.3 | U | 8.1 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.80 | U | 8.1 | 0.80 |
| 132-64-9 | Dibenzofuran | 0.80 | U | 61 | 0.80 |
| 84-66-2 | Diethyl phthalate | 19 | U | 85 | 19 |
| 131-11-3 | Dimethyl phthalate | 21 | U | 85 | 21 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 85 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.6 | U | 85 | 9.6 |
| 206-44-0 | Fluoranthene | 0.67 | U | 8.1 | 0.67 |
| 86-73-7 | Fluorene | 0.64 | U | 8.1 | 0.64 |
| 118-74-1 | Hexachlorobenzene | 2.6 | U | 8.1 | 2.6 |
| 87-68-3 | Hexachlorobutadiene | 6.8 | U | 61 | 6.8 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.8 | U | 400 | 9.8 |
| 67-72-1 | Hexachloroethane | 11 | U | 61 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.43 | U | 8.1 | 0.43 |
| 78-59-1 | Isophorone | 16 | U | 61 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.7 | U | 61 | 7.7 |
| 86-30-6 | N-Nitrosodiphenylamine | 26 | U | 61 | 26 |
| 91-20-3 | Naphthalene | 1.0 | U | 8.1 | 1.0 |
| 98-95-3 | Nitrobenzene | 2.7 | U | 120 | 2.7 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.89 | U | 8.1 | 0.89 |
| 108-95-2 | Phenol | 8.9 | U | 61 | 8.9 |
| 129-00-0 | Pyrene | 0.53 | U | 8.1 | 0.53 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 490 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 51111008.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:04

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.03(g) | Date Analyzed: 11/11/2015 12:05

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 17.8 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 73 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 60 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 58 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 61 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 63 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 41 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 51111009.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:06

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.84(g) | Date Analyzed: 11/11/2015 12:30

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 60 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.54 | U | 60 | 0.54 |
| 95-57-8 | 2-Chlorophenol | 9.8 | U | 60 | 9.8 |
| 91-57-6 | 2-Methylnaphthalene | 0.60 | U | 7.9 | 0.60 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.9 | U | 60 | 9.9 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 60 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 60 | 15 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.91 | U | 7.9 | 0.91 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 7.9 | 0.42 |
| 98-86-2 | Acetophenone | 13 | J | 120 | 11 |
| 120-12-7 | Anthracene | 0.93 | U | 7.9 | 0.93 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.75 | U | 7.9 | 0.75 |
| 50-32-8 | Benzo[a]pyrene | 0.76 | U | 7.9 | 0.76 |
| 205-99-2 | Benzo[b]fluoranthene | 0.70 | U | 7.9 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.42 | U | 7.9 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 51111009.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:06

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.84(g) | Date Analyzed: 11/11/2015 12:30

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.81 | U | 7.9 | 0.81 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 33 | J | 83 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 83 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 60 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 7.9 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.79 | U | 7.9 | 0.79 |
| 132-64-9 | Dibenzofuran | 0.79 | U | 60 | 0.79 |
| 84-66-2 | Diethyl phthalate | 19 | U | 83 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 83 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 83 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.4 | U | 83 | 9.4 |
| 206-44-0 | Fluoranthene | 0.66 | U | 7.9 | 0.66 |
| 86-73-7 | Fluorene | 0.63 | U | 7.9 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.9 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.7 | U | 60 | 6.7 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.7 | U | 390 | 9.7 |
| 67-72-1 | Hexachloroethane | 11 | U | 60 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.42 | U | 7.9 | 0.42 |
| 78-59-1 | Isophorone | 15 | U | 60 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.5 | U | 60 | 7.5 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 60 | 25 |
| 91-20-3 | Naphthalene | 0.98 | U | 7.9 | 0.98 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.87 | U | 7.9 | 0.87 |
| 108-95-2 | Phenol | 8.7 | U | 60 | 8.7 |
| 129-00-0 | Pyrene | 0.52 | U | 7.9 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 51111009.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:06

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.84(g) | Date Analyzed: 11/11/2015 12:30

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 85 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 72 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 69 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 72 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 72 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 42 | | 10-110 |

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Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 51111010.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:08

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.28(g) | Date Analyzed: 11/11/2015 12:54

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 58 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 29 | U | 180 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10 | U | 180 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 23 | U | 180 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 23 | U | 180 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 230 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.53 | U | 58 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 9.6 | U | 58 | 9.6 |
| 91-57-6 | 2-Methylnaphthalene | 0.58 | U | 7.8 | 0.58 |
| 95-48-7 | 2-Methylphenol | 13 | U | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.7 | U | 58 | 9.7 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 58 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 58 | 15 |
| 100-01-6 | 4-Nitroaniline | 30 | U | 230 | 30 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.89 | U | 7.8 | 0.89 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.8 | 0.41 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.91 | U | 7.8 | 0.91 |
| 1912-24-9 | Atrazine | 11 | U | 230 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.74 | U | 7.8 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 0.75 | U | 7.8 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 0.69 | U | 7.8 | 0.69 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.41 | U | 7.8 | 0.41 |

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GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 51111010.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:08

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.28(g) | Date Analyzed: 11/11/2015 12:54

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.79 | U | 7.8 | 0.79 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.3 | U | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 38 | J | 82 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 82 | 12 |
| 105-60-2 | Caprolactam | 43 | U | 390 | 43 |
| 86-74-8 | Carbazole | 32 | U | 58 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 7.8 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.77 | U | 7.8 | 0.77 |
| 132-64-9 | Dibenzofuran | 0.77 | U | 58 | 0.77 |
| 84-66-2 | Diethyl phthalate | 19 | U | 82 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 82 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 82 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.2 | U | 82 | 9.2 |
| 206-44-0 | Fluoranthene | 0.64 | U | 7.8 | 0.64 |
| 86-73-7 | Fluorene | 0.62 | U | 7.8 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.8 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.5 | U | 58 | 6.5 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.5 | U | 390 | 9.5 |
| 67-72-1 | Hexachloroethane | 11 | U | 58 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.41 | U | 7.8 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 58 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.4 | U | 58 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 58 | 25 |
| 91-20-3 | Naphthalene | 0.96 | U | 7.8 | 0.96 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.85 | U | 7.8 | 0.85 |
| 108-95-2 | Phenol | 8.5 | U | 58 | 8.5 |
| 129-00-0 | Pyrene | 0.51 | U | 7.8 | 0.51 |
| 15831-10-4 | 3 & 4 Methylphenol | 23 | U | 470 | 23 |

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GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 51111010.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:08

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.28(g) | Date Analyzed: 11/11/2015 12:54

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 77 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 67 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 64 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 68 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 69 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 25 | | 10-110 |

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GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 51111011.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:10

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.15(g) | Date Analyzed: 11/11/2015 13:19

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 60 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.54 | U | 60 | 0.54 |
| 95-57-8 | 2-Chlorophenol | 9.8 | U | 60 | 9.8 |
| 91-57-6 | 2-Methylnaphthalene | 0.60 | U | 8.0 | 0.60 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.9 | U | 60 | 9.9 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 60 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 60 | 15 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.91 | U | 8.0 | 0.91 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 8.0 | 0.42 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.93 | U | 8.0 | 0.93 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.75 | U | 8.0 | 0.75 |
| 50-32-8 | Benzo[a]pyrene | 0.76 | U | 8.0 | 0.76 |
| 205-99-2 | Benzo[b]fluoranthene | 0.70 | U | 8.0 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.42 | U | 8.0 | 0.42 |

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GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 51111011.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:10

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.15(g) | Date Analyzed: 11/11/2015 13:19

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.81 | U | 8.0 | 0.81 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 32 | J | 83 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 83 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 60 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 8.0 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.79 | U | 8.0 | 0.79 |
| 132-64-9 | Dibenzofuran | 0.79 | U | 60 | 0.79 |
| 84-66-2 | Diethyl phthalate | 19 | U | 83 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 83 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 83 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.4 | U | 83 | 9.4 |
| 206-44-0 | Fluoranthene | 0.66 | U | 8.0 | 0.66 |
| 86-73-7 | Fluorene | 0.63 | U | 8.0 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 8.0 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.7 | U | 60 | 6.7 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.7 | U | 390 | 9.7 |
| 67-72-1 | Hexachloroethane | 11 | U | 60 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.42 | U | 8.0 | 0.42 |
| 78-59-1 | Isophorone | 15 | U | 60 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.5 | U | 60 | 7.5 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 60 | 25 |
| 91-20-3 | Naphthalene | 0.98 | U | 8.0 | 0.98 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.87 | U | 8.0 | 0.87 |
| 108-95-2 | Phenol | 8.7 | U | 60 | 8.7 |
| 129-00-0 | Pyrene | 0.52 | U | 8.0 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

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GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 51111011.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:10

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.15(g) | Date Analyzed: 11/11/2015 13:19

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 79 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 60 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 62 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 62 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 68 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 33 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 51111012.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:12

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.53(g) | Date Analyzed: 11/11/2015 13:44

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 61 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 12 | U | 120 | 12 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 400 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 21 | U | 240 | 21 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.55 | U | 61 | 0.55 |
| 95-57-8 | 2-Chlorophenol | 9.9 | U | 61 | 9.9 |
| 91-57-6 | 2-Methylnaphthalene | 0.61 | U | 8.1 | 0.61 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 10 | U | 61 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 61 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 21 | U | 180 | 21 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 61 | 16 |
| 100-01-6 | 4-Nitroaniline | 32 | U | 240 | 32 |
| 100-02-7 | 4-Nitrophenol | 21 | U | 400 | 21 |
| 83-32-9 | Acenaphthene | 0.92 | U | 8.1 | 0.92 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 8.1 | 0.42 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.95 | U | 8.1 | 0.95 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 15 | U | 120 | 15 |
| 56-55-3 | Benzo[a]anthracene | 0.76 | U | 8.1 | 0.76 |
| 50-32-8 | Benzo[a]pyrene | 0.78 | U | 8.1 | 0.78 |
| 205-99-2 | Benzo[b]fluoranthene | 0.72 | U | 8.1 | 0.72 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.42 | U | 8.1 | 0.42 |

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GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 51111012.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:12

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.53(g) | Date Analyzed: 11/11/2015 13:44

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.82 | U | 8.1 | 0.82 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 27 | U | 120 | 27 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 32 | J | 85 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 85 | 12 |
| 105-60-2 | Caprolactam | 45 | U | 400 | 45 |
| 86-74-8 | Carbazole | 33 | U | 61 | 33 |
| 218-01-9 | Chrysene | 1.3 | U | 8.1 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.80 | U | 8.1 | 0.80 |
| 132-64-9 | Dibenzofuran | 0.80 | U | 61 | 0.80 |
| 84-66-2 | Diethyl phthalate | 19 | U | 85 | 19 |
| 131-11-3 | Dimethyl phthalate | 21 | U | 85 | 21 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 85 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.6 | U | 85 | 9.6 |
| 206-44-0 | Fluoranthene | 0.67 | U | 8.1 | 0.67 |
| 86-73-7 | Fluorene | 0.64 | U | 8.1 | 0.64 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 8.1 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.8 | U | 61 | 6.8 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.8 | U | 400 | 9.8 |
| 67-72-1 | Hexachloroethane | 11 | U | 61 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.42 | U | 8.1 | 0.42 |
| 78-59-1 | Isophorone | 16 | U | 61 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.6 | U | 61 | 7.6 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 61 | 25 |
| 91-20-3 | Naphthalene | 0.99 | U | 8.1 | 0.99 |
| 98-95-3 | Nitrobenzene | 2.7 | U | 120 | 2.7 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.89 | U | 8.1 | 0.89 |
| 108-95-2 | Phenol | 8.9 | U | 61 | 8.9 |
| 129-00-0 | Pyrene | 0.53 | U | 8.1 | 0.53 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 490 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 51111012.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:12

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.53(g) | Date Analyzed: 11/11/2015 13:44

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 80 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 64 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 62 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 67 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 70 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 38 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 51111013.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:14

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.20(g) | Date Analyzed: 11/11/2015 14:08

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 61 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 12 | U | 120 | 12 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 400 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 21 | U | 240 | 21 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.55 | U | 61 | 0.55 |
| 95-57-8 | 2-Chlorophenol | 9.9 | U | 61 | 9.9 |
| 91-57-6 | 2-Methylnaphthalene | 0.61 | U | 8.1 | 0.61 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 10 | U | 61 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 61 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 21 | U | 180 | 21 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 61 | 16 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 21 | U | 400 | 21 |
| 83-32-9 | Acenaphthene | 0.92 | U | 8.1 | 0.92 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 8.1 | 0.42 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.94 | U | 8.1 | 0.94 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 15 | U | 120 | 15 |
| 56-55-3 | Benzo[a]anthracene | 0.76 | U | 8.1 | 0.76 |
| 50-32-8 | Benzo[a]pyrene | 0.78 | U | 8.1 | 0.78 |
| 205-99-2 | Benzo[b]fluoranthene | 0.71 | U | 8.1 | 0.71 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.42 | U | 8.1 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 51111013.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:14

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.20(g) | Date Analyzed: 11/11/2015 14:08

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.82 | U | 8.1 | 0.82 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 27 | U | 120 | 27 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 23 | U | 85 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 85 | 12 |
| 105-60-2 | Caprolactam | 45 | U | 400 | 45 |
| 86-74-8 | Carbazole | 33 | U | 61 | 33 |
| 218-01-9 | Chrysene | 1.3 | U | 8.1 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.80 | U | 8.1 | 0.80 |
| 132-64-9 | Dibenzofuran | 0.80 | U | 61 | 0.80 |
| 84-66-2 | Diethyl phthalate | 19 | U | 85 | 19 |
| 131-11-3 | Dimethyl phthalate | 21 | U | 85 | 21 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 85 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.6 | U | 85 | 9.6 |
| 206-44-0 | Fluoranthene | 0.67 | U | 8.1 | 0.67 |
| 86-73-7 | Fluorene | 0.64 | U | 8.1 | 0.64 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 8.1 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.8 | U | 61 | 6.8 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.8 | U | 400 | 9.8 |
| 67-72-1 | Hexachloroethane | 11 | U | 61 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.42 | U | 8.1 | 0.42 |
| 78-59-1 | Isophorone | 16 | U | 61 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.6 | U | 61 | 7.6 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 61 | 25 |
| 91-20-3 | Naphthalene | 0.99 | U | 8.1 | 0.99 |
| 98-95-3 | Nitrobenzene | 2.7 | U | 120 | 2.7 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.88 | U | 8.1 | 0.88 |
| 108-95-2 | Phenol | 8.8 | U | 61 | 8.8 |
| 129-00-0 | Pyrene | 0.53 | U | 8.1 | 0.53 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 51111013.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:14

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30.20(g) | Date Analyzed: 11/11/2015 14:08

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 70 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 54 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 52 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 56 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 56 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 32 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 51111014.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:16

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.67(g) | Date Analyzed: 11/11/2015 14:33

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 19.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.4 | U | 62 | 4.4 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 12 | U | 120 | 12 |
| 95-95-4 | 2,4,5-Trichlorophenol | 31 | U | 190 | 31 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 190 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 25 | U | 190 | 25 |
| 105-67-9 | 2,4-Dimethylphenol | 25 | U | 190 | 25 |
| 51-28-5 | 2,4-Dinitrophenol | 26 | U | 410 | 26 |
| 121-14-2 | 2,4-Dinitrotoluene | 21 | U | 250 | 21 |
| 606-20-2 | 2,6-Dinitrotoluene | 26 | U | 250 | 26 |
| 91-58-7 | 2-Chloronaphthalene | 0.56 | U | 62 | 0.56 |
| 95-57-8 | 2-Chlorophenol | 10 | U | 62 | 10 |
| 91-57-6 | 2-Methylnaphthalene | 0.62 | U | 8.3 | 0.62 |
| 95-48-7 | 2-Methylphenol | 14 | U | 250 | 14 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 250 | 11 |
| 88-75-5 | 2-Nitrophenol | 10 | U | 62 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 20 | U | 250 | 20 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 190 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 62 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 26 | U | 190 | 26 |
| 106-47-8 | 4-Chloroaniline | 21 | U | 190 | 21 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 62 | 16 |
| 100-01-6 | 4-Nitroaniline | 32 | U | 250 | 32 |
| 100-02-7 | 4-Nitrophenol | 21 | U | 410 | 21 |
| 83-32-9 | Acenaphthene | 0.95 | U | 8.3 | 0.95 |
| 208-96-8 | Acenaphthylene | 0.44 | U | 8.3 | 0.44 |
| 98-86-2 | Acetophenone | 13 | J | 120 | 11 |
| 120-12-7 | Anthracene | 0.97 | U | 8.3 | 0.97 |
| 1912-24-9 | Atrazine | 11 | U | 250 | 11 |
| 100-52-7 | Benzaldehyde | 15 | U | 120 | 15 |
| 56-55-3 | Benzo[a]anthracene | 0.79 | U | 8.3 | 0.79 |
| 50-32-8 | Benzo[a]pyrene | 0.80 | U | 8.3 | 0.80 |
| 205-99-2 | Benzo[b]fluoranthene | 0.74 | U | 8.3 | 0.74 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.44 | U | 8.3 | 0.44 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 51111014.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:16

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.67(g) | Date Analyzed: 11/11/2015 14:33

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 19.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.85 | U | 8.3 | 0.85 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 27 | U | 120 | 27 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.5 | U | 120 | 2.5 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 40 | J | 87 | 24 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 87 | 12 |
| 105-60-2 | Caprolactam | 46 | U | 410 | 46 |
| 86-74-8 | Carbazole | 34 | U | 62 | 34 |
| 218-01-9 | Chrysene | 1.4 | U | 8.3 | 1.4 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.82 | U | 8.3 | 0.82 |
| 132-64-9 | Dibenzofuran | 0.82 | U | 62 | 0.82 |
| 84-66-2 | Diethyl phthalate | 20 | U | 87 | 20 |
| 131-11-3 | Dimethyl phthalate | 21 | U | 87 | 21 |
| 84-74-2 | Di-n-butyl phthalate | 19 | U | 87 | 19 |
| 117-84-0 | Di-n-octyl phthalate | 9.9 | U | 87 | 9.9 |
| 206-44-0 | Fluoranthene | 0.69 | U | 8.3 | 0.69 |
| 86-73-7 | Fluorene | 0.66 | U | 8.3 | 0.66 |
| 118-74-1 | Hexachlorobenzene | 2.6 | U | 8.3 | 2.6 |
| 87-68-3 | Hexachlorobutadiene | 7.0 | U | 62 | 7.0 |
| 77-47-4 | Hexachlorocyclopentadiene | 10 | U | 410 | 10 |
| 67-72-1 | Hexachloroethane | 11 | U | 62 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.44 | U | 8.3 | 0.44 |
| 78-59-1 | Isophorone | 16 | U | 62 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.9 | U | 62 | 7.9 |
| 86-30-6 | N-Nitrosodiphenylamine | 26 | U | 62 | 26 |
| 91-20-3 | Naphthalene | 1.0 | U | 8.3 | 1.0 |
| 98-95-3 | Nitrobenzene | 2.7 | U | 120 | 2.7 |
| 87-86-5 | Pentachlorophenol | 11 | U | 190 | 11 |
| 85-01-8 | Phenanthrene | 0.91 | U | 8.3 | 0.91 |
| 108-95-2 | Phenol | 9.1 | U | 62 | 9.1 |
| 129-00-0 | Pyrene | 0.55 | U | 8.3 | 0.55 |
| 15831-10-4 | 3 & 4 Methylphenol | 25 | U | 500 | 25 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 51111014.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:16

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.67(g) | Date Analyzed: 11/11/2015 14:33

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 19.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 87 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 68 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 68 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 71 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 73 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 45 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 51111015.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:18

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.78(g) | Date Analyzed: 11/11/2015 14:57

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 13.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 58 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 29 | U | 170 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10 | U | 170 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 23 | U | 170 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 23 | U | 170 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 24 | U | 380 | 24 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 24 | U | 230 | 24 |
| 91-58-7 | 2-Chloronaphthalene | 0.52 | U | 58 | 0.52 |
| 95-57-8 | 2-Chlorophenol | 9.5 | U | 58 | 9.5 |
| 91-57-6 | 2-Methylnaphthalene | 0.58 | U | 7.8 | 0.58 |
| 95-48-7 | 2-Methylphenol | 13 | U | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.7 | U | 58 | 9.7 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 170 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 58 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 24 | U | 170 | 24 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 170 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 58 | 15 |
| 100-01-6 | 4-Nitroaniline | 30 | U | 230 | 30 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 380 | 20 |
| 83-32-9 | Acenaphthene | 0.88 | U | 7.8 | 0.88 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.8 | 0.41 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.91 | U | 7.8 | 0.91 |
| 1912-24-9 | Atrazine | 11 | U | 230 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.73 | U | 7.8 | 0.73 |
| 50-32-8 | Benzo[a]pyrene | 0.74 | U | 7.8 | 0.74 |
| 205-99-2 | Benzo[b]fluoranthene | 0.69 | U | 7.8 | 0.69 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.41 | U | 7.8 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 51111015.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:18

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.78(g) | Date Analyzed: 11/11/2015 14:57

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 13.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.79 | U | 7.8 | 0.79 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.3 | U | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 29 | J | 81 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 81 | 12 |
| 105-60-2 | Caprolactam | 43 | U | 380 | 43 |
| 86-74-8 | Carbazole | 31 | U | 58 | 31 |
| 218-01-9 | Chrysene | 1.3 | U | 7.8 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.77 | U | 7.8 | 0.77 |
| 132-64-9 | Dibenzofuran | 0.77 | U | 58 | 0.77 |
| 84-66-2 | Diethyl phthalate | 19 | U | 81 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 81 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 17 | U | 81 | 17 |
| 117-84-0 | Di-n-octyl phthalate | 9.2 | U | 81 | 9.2 |
| 206-44-0 | Fluoranthene | 0.64 | U | 7.8 | 0.64 |
| 86-73-7 | Fluorene | 0.62 | U | 7.8 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 2.4 | U | 7.8 | 2.4 |
| 87-68-3 | Hexachlorobutadiene | 6.5 | U | 58 | 6.5 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.4 | U | 380 | 9.4 |
| 67-72-1 | Hexachloroethane | 10 | U | 58 | 10 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.41 | U | 7.8 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 58 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.3 | U | 58 | 7.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 24 | U | 58 | 24 |
| 91-20-3 | Naphthalene | 0.95 | U | 7.8 | 0.95 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 170 | 11 |
| 85-01-8 | Phenanthrene | 0.85 | U | 7.8 | 0.85 |
| 108-95-2 | Phenol | 8.5 | U | 58 | 8.5 |
| 129-00-0 | Pyrene | 0.51 | U | 7.8 | 0.51 |
| 15831-10-4 | 3 & 4 Methylphenol | 23 | U | 470 | 23 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 51111015.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:18

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 29.78(g) | Date Analyzed: 11/11/2015 14:57

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 13.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 78 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 61 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 60 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 59 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 64 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 40 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 51111019.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:20

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.86(g) | Date Analyzed: 11/11/2015 14:50

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 60 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 400 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.54 | U | 60 | 0.54 |
| 95-57-8 | 2-Chlorophenol | 9.9 | U | 60 | 9.9 |
| 91-57-6 | 2-Methylnaphthalene | 0.60 | U | 8.0 | 0.60 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 10 | U | 60 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 60 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 60 | 16 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 400 | 20 |
| 83-32-9 | Acenaphthene | 0.91 | U | 8.0 | 0.91 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 8.0 | 0.42 |
| 98-86-2 | Acetophenone | 38 | J | 120 | 11 |
| 120-12-7 | Anthracene | 0.94 | U | 8.0 | 0.94 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 19 | J | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.76 | U | 8.0 | 0.76 |
| 50-32-8 | Benzo[a]pyrene | 0.77 | U | 8.0 | 0.77 |
| 205-99-2 | Benzo[b]fluoranthene | 0.71 | U | 8.0 | 0.71 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.42 | U | 8.0 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 51111019.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:20

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.86(g) | Date Analyzed: 11/11/2015 14:50

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.82 | U | 8.0 | 0.82 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 23 | U | 84 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 84 | 12 |
| 105-60-2 | Caprolactam | 45 | U | 400 | 45 |
| 86-74-8 | Carbazole | 32 | U | 60 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 8.0 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.79 | U | 8.0 | 0.79 |
| 132-64-9 | Dibenzofuran | 0.79 | U | 60 | 0.79 |
| 84-66-2 | Diethyl phthalate | 19 | U | 84 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 84 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 84 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.5 | U | 84 | 9.5 |
| 206-44-0 | Fluoranthene | 0.66 | U | 8.0 | 0.66 |
| 86-73-7 | Fluorene | 0.64 | U | 8.0 | 0.64 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 8.0 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.7 | U | 60 | 6.7 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.7 | U | 400 | 9.7 |
| 67-72-1 | Hexachloroethane | 11 | U | 60 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.42 | U | 8.0 | 0.42 |
| 78-59-1 | Isophorone | 16 | U | 60 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.6 | U | 60 | 7.6 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 60 | 25 |
| 91-20-3 | Naphthalene | 0.99 | U | 8.0 | 0.99 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.88 | U | 8.0 | 0.88 |
| 108-95-2 | Phenol | 8.8 | U | 60 | 8.8 |
| 129-00-0 | Pyrene | 0.53 | U | 8.0 | 0.53 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 51111019.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:20

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.86(g) | Date Analyzed: 11/11/2015 14:50

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 88 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 80 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 80 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 78 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 73 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 40 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 51111020.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:22

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.84(g) | Date Analyzed: 11/11/2015 15:14

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 60 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.54 | U | 60 | 0.54 |
| 95-57-8 | 2-Chlorophenol | 9.8 | U | 60 | 9.8 |
| 91-57-6 | 2-Methylnaphthalene | 0.60 | U | 8.0 | 0.60 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.9 | U | 60 | 9.9 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 60 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 60 | 16 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.91 | U | 8.0 | 0.91 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 8.0 | 0.42 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.93 | U | 8.0 | 0.93 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.75 | U | 8.0 | 0.75 |
| 50-32-8 | Benzo[a]pyrene | 0.77 | U | 8.0 | 0.77 |
| 205-99-2 | Benzo[b]fluoranthene | 0.71 | U | 8.0 | 0.71 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.42 | U | 8.0 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 51111020.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:22

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.84(g) | Date Analyzed: 11/11/2015 15:14

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.81 | U | 8.0 | 0.81 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 23 | U | 84 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 84 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 60 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 8.0 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.79 | U | 8.0 | 0.79 |
| 132-64-9 | Dibenzofuran | 0.79 | U | 60 | 0.79 |
| 84-66-2 | Diethyl phthalate | 19 | U | 84 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 84 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 84 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.4 | U | 84 | 9.4 |
| 206-44-0 | Fluoranthene | 0.66 | U | 8.0 | 0.66 |
| 86-73-7 | Fluorene | 0.63 | U | 8.0 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 8.0 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.7 | U | 60 | 6.7 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.7 | U | 390 | 9.7 |
| 67-72-1 | Hexachloroethane | 11 | U | 60 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.42 | U | 8.0 | 0.42 |
| 78-59-1 | Isophorone | 16 | U | 60 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.5 | U | 60 | 7.5 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 60 | 25 |
| 91-20-3 | Naphthalene | 0.98 | U | 8.0 | 0.98 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.87 | U | 8.0 | 0.87 |
| 108-95-2 | Phenol | 8.7 | U | 60 | 8.7 |
| 129-00-0 | Pyrene | 0.53 | U | 8.0 | 0.53 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 51111020.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:22

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.84(g) | Date Analyzed: 11/11/2015 15:14

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 72 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 64 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 67 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 67 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 63 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 32 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 51111021.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:24

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30.37(g) | Date Analyzed: 11/11/2015 15:39

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 58 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 29 | U | 170 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10 | U | 170 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 23 | U | 170 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 23 | U | 170 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 24 | U | 380 | 24 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 24 | U | 230 | 24 |
| 91-58-7 | 2-Chloronaphthalene | 0.52 | U | 58 | 0.52 |
| 95-57-8 | 2-Chlorophenol | 9.5 | U | 58 | 9.5 |
| 91-57-6 | 2-Methylnaphthalene | 0.58 | U | 7.7 | 0.58 |
| 95-48-7 | 2-Methylphenol | 13 | U | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.6 | U | 58 | 9.6 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 170 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 58 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 24 | U | 170 | 24 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 170 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 58 | 15 |
| 100-01-6 | 4-Nitroaniline | 30 | U | 230 | 30 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 380 | 20 |
| 83-32-9 | Acenaphthene | 0.88 | U | 7.7 | 0.88 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.7 | 0.41 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.90 | U | 7.7 | 0.90 |
| 1912-24-9 | Atrazine | 11 | U | 230 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.73 | U | 7.7 | 0.73 |
| 50-32-8 | Benzo[a]pyrene | 0.74 | U | 7.7 | 0.74 |
| 205-99-2 | Benzo[b]fluoranthene | 0.68 | U | 7.7 | 0.68 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.41 | U | 7.7 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 51111021.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:24

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30.37(g) | Date Analyzed: 11/11/2015 15:39

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.79 | U | 7.7 | 0.79 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 25 | U | 120 | 25 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.3 | U | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 60 | J | 81 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 81 | 12 |
| 105-60-2 | Caprolactam | 43 | U | 380 | 43 |
| 86-74-8 | Carbazole | 31 | U | 58 | 31 |
| 218-01-9 | Chrysene | 1.3 | U | 7.7 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.76 | U | 7.7 | 0.76 |
| 132-64-9 | Dibenzofuran | 0.76 | U | 58 | 0.76 |
| 84-66-2 | Diethyl phthalate | 19 | U | 81 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 81 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | J | 81 | 17 |
| 117-84-0 | Di-n-octyl phthalate | 9.1 | U | 81 | 9.1 |
| 206-44-0 | Fluoranthene | 0.64 | U | 7.7 | 0.64 |
| 86-73-7 | Fluorene | 0.61 | U | 7.7 | 0.61 |
| 118-74-1 | Hexachlorobenzene | 2.4 | U | 7.7 | 2.4 |
| 87-68-3 | Hexachlorobutadiene | 6.5 | U | 58 | 6.5 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.4 | U | 380 | 9.4 |
| 67-72-1 | Hexachloroethane | 10 | U | 58 | 10 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.41 | U | 7.7 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 58 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.3 | U | 58 | 7.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 24 | U | 58 | 24 |
| 91-20-3 | Naphthalene | 0.95 | U | 7.7 | 0.95 |
| 98-95-3 | Nitrobenzene | 2.5 | U | 120 | 2.5 |
| 87-86-5 | Pentachlorophenol | 11 | U | 170 | 11 |
| 85-01-8 | Phenanthrene | 0.85 | U | 7.7 | 0.85 |
| 108-95-2 | Phenol | 8.5 | U | 58 | 8.5 |
| 129-00-0 | Pyrene | 0.51 | U | 7.7 | 0.51 |
| 15831-10-4 | 3 & 4 Methylphenol | 23 | U | 460 | 23 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 51111021.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:24

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30.37(g) | Date Analyzed: 11/11/2015 15:39

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 77 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 57 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 65 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 65 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 59 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 33 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 51111022.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:26

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.64(g) | Date Analyzed: 11/11/2015 16:03

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 59 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.53 | U | 59 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 9.7 | U | 59 | 9.7 |
| 91-57-6 | 2-Methylnaphthalene | 0.59 | U | 7.9 | 0.59 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.8 | U | 59 | 9.8 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 59 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 59 | 15 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.90 | U | 7.9 | 0.90 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.9 | 0.41 |
| 98-86-2 | Acetophenone | 11 | J | 120 | 11 |
| 120-12-7 | Anthracene | 0.92 | U | 7.9 | 0.92 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.75 | U | 7.9 | 0.75 |
| 50-32-8 | Benzo[a]pyrene | 0.76 | U | 7.9 | 0.76 |
| 205-99-2 | Benzo[b]fluoranthene | 0.70 | U | 7.9 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.41 | U | 7.9 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 51111022.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:26

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.64(g) | Date Analyzed: 11/11/2015 16:03

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.80 | U | 7.9 | 0.80 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 22 | U | 83 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 83 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 59 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 7.9 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.78 | U | 7.9 | 0.78 |
| 132-64-9 | Dibenzofuran | 0.78 | U | 59 | 0.78 |
| 84-66-2 | Diethyl phthalate | 19 | U | 83 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 83 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 83 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.4 | U | 83 | 9.4 |
| 206-44-0 | Fluoranthene | 0.65 | U | 7.9 | 0.65 |
| 86-73-7 | Fluorene | 0.63 | U | 7.9 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.9 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.6 | U | 59 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.6 | U | 390 | 9.6 |
| 67-72-1 | Hexachloroethane | 11 | U | 59 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.41 | U | 7.9 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 59 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.5 | U | 59 | 7.5 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 59 | 25 |
| 91-20-3 | Naphthalene | 0.97 | U | 7.9 | 0.97 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.86 | U | 7.9 | 0.86 |
| 108-95-2 | Phenol | 8.6 | U | 59 | 8.6 |
| 129-00-0 | Pyrene | 0.52 | U | 7.9 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 470 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 51111022.D

Analysis Method: 8270D | Date Collected: 11/05/2015 09:26

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.64(g) | Date Analyzed: 11/11/2015 16:03

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 79 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 62 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 73 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 70 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 68 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 28 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 51111023.D

Analysis Method: 8270D | Date Collected: 11/05/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30.09(g) | Date Analyzed: 11/11/2015 16:28

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 58 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 29 | U | 170 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10 | U | 170 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 23 | U | 170 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 23 | U | 170 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 24 | U | 380 | 24 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 24 | U | 230 | 24 |
| 91-58-7 | 2-Chloronaphthalene | 0.52 | U | 58 | 0.52 |
| 95-57-8 | 2-Chlorophenol | 9.6 | U | 58 | 9.6 |
| 91-57-6 | 2-Methylnaphthalene | 0.58 | U | 7.8 | 0.58 |
| 95-48-7 | 2-Methylphenol | 13 | U | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.7 | U | 58 | 9.7 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 170 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 58 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 24 | U | 170 | 24 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 170 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 58 | 15 |
| 100-01-6 | 4-Nitroaniline | 30 | U | 230 | 30 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 380 | 20 |
| 83-32-9 | Acenaphthene | 0.89 | U | 7.8 | 0.89 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.8 | 0.41 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.91 | U | 7.8 | 0.91 |
| 1912-24-9 | Atrazine | 11 | U | 230 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.73 | U | 7.8 | 0.73 |
| 50-32-8 | Benzo[a]pyrene | 0.75 | U | 7.8 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 0.69 | U | 7.8 | 0.69 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.41 | U | 7.8 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 51111023.D

Analysis Method: 8270D | Date Collected: 11/05/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30.09(g) | Date Analyzed: 11/11/2015 16:28

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.79 | U | 7.8 | 0.79 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.3 | U | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 22 | J | 82 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 82 | 12 |
| 105-60-2 | Caprolactam | 43 | U | 380 | 43 |
| 86-74-8 | Carbazole | 31 | U | 58 | 31 |
| 218-01-9 | Chrysene | 1.3 | U | 7.8 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.77 | U | 7.8 | 0.77 |
| 132-64-9 | Dibenzofuran | 0.77 | U | 58 | 0.77 |
| 84-66-2 | Diethyl phthalate | 19 | U | 82 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 82 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 17 | U | 82 | 17 |
| 117-84-0 | Di-n-octyl phthalate | 9.2 | U | 82 | 9.2 |
| 206-44-0 | Fluoranthene | 0.64 | U | 7.8 | 0.64 |
| 86-73-7 | Fluorene | 0.62 | U | 7.8 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 2.4 | U | 7.8 | 2.4 |
| 87-68-3 | Hexachlorobutadiene | 6.5 | U | 58 | 6.5 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.4 | U | 380 | 9.4 |
| 67-72-1 | Hexachloroethane | 10 | U | 58 | 10 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.41 | U | 7.8 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 58 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.3 | U | 58 | 7.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 24 | U | 58 | 24 |
| 91-20-3 | Naphthalene | 0.96 | U | 7.8 | 0.96 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 170 | 11 |
| 85-01-8 | Phenanthrene | 0.85 | U | 7.8 | 0.85 |
| 108-95-2 | Phenol | 8.5 | U | 58 | 8.5 |
| 129-00-0 | Pyrene | 0.51 | U | 7.8 | 0.51 |
| 15831-10-4 | 3 & 4 Methylphenol | 23 | U | 470 | 23 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 51111023.D

Analysis Method: 8270D | Date Collected: 11/05/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30.09(g) | Date Analyzed: 11/11/2015 16:28

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 81 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 60 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 74 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 71 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 68 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 24 | | 10-110 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48  Calibration End Date: 10/02/2015 14:04  Calibration ID: 30980

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-200091/6 | 51002006.D |
| Level 2 | STD2 240-200091/5 | 51002005.D |
| Level 3 | STD3 240-200091/4 | 51002004.D |
| Level 4 | STD4 240-200091/3 | 51002003.D |
| Level 5 | STD5 240-200091/2 | 51002002.D |
| Level 6 | STD6 240-200091/10 | 51002010.D |
| Level 7 | STD7 240-200091/9 | 51002009.D |
| Level 8 | STD8 240-200091/8 | 51002008.D |
| Level 9 | STD9 240-200091/7 | 51002007.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,4-Dioxane | 0.6695 | 0.7664<br>0.6450 | 0.7112<br>0.6601 | 0.6791<br>0.6797 | 0.7113 | Ave | | 0.6903 | | | | 5.6 | | 20.0 | | | |
| N-Nitrosodimethylamine | 0.9213 | 0.9861<br>0.9036 | 0.9021<br>0.9073 | 0.9106<br>0.9353 | 0.9364 | Ave | | 0.9253 | | | | 3.0 | | 20.0 | | | |
| Pyridine | 1.6752 | 1.7590<br>1.6375 | 1.6236<br>1.6765 | 1.6397<br>1.7157 | 1.7460 | Ave | | 1.6842 | | | | 3.0 | | 20.0 | | | |
| Benzaldehyde | 1.1161 | 1.2599<br>0.9941 | 1.2278<br>0.9643 | 1.2174<br>0.9002 | 1.2082 | Ave | | 1.1110 | | | 0.0100 | 12.6 | | 20.0 | | | |
| Phenol | 1.8772 | 2.0486<br>1.7989 | 1.9063<br>1.8076 | 1.9409<br>1.8084 | 1.9432 | Ave | | 1.8914 | | | 0.8000 | 4.6 | | 20.0 | | | |
| Aniline | 2.4273 | 2.4891<br>2.3113 | 2.4030<br>2.3297 | 2.4066<br>2.3780 | 2.4676 | Ave | | 2.4016 | | | | 2.6 | | 20.0 | | | |
| Bis(2-chloroethyl)ether | 1.5242 | 1.5993<br>1.4438 | 1.5227<br>1.4462 | 1.5177<br>1.4436 | 1.5534 | Ave | | 1.5064 | | | 0.7000 | 3.8 | | 20.0 | | | |
| 2-Chlorophenol | 1.5436 | 1.5843<br>1.4930 | 1.5903<br>1.5124 | 1.5578<br>1.5091 | 1.5573 | Ave | | 1.5435 | | | 0.8000 | 2.3 | | 20.0 | | | |
| n-Decane | 1.8806 | 2.1172<br>1.7823 | 2.0495<br>1.7515 | 1.9665<br>1.7279 | 2.0152 | Ave | | 1.9113 | | | | 7.7 | | 20.0 | | | |
| 1,3-Dichlorobenzene | 1.5929 | 1.6788<br>1.5009 | 1.5686<br>1.5104 | 1.6231<br>1.5138 | 1.6452 | Ave | | 1.5792 | | | | 4.3 | | 20.0 | | | |
| 1,4-Dichlorobenzene | 1.6286 | 1.7289<br>1.5274 | 1.6483<br>1.5572 | 1.6056<br>1.5444 | 1.6767 | Ave | | 1.6146 | | | | 4.3 | | 20.0 | | | |
| Benzyl alcohol | 1.0249 | 1.1085<br>0.9791 | 1.0175<br>0.9932 | 1.0225<br>0.9932 | 1.0587 | Ave | | 1.0247 | | | | 4.1 | | 20.0 | | | |
| 1,2-Dichlorobenzene | 1.5382 | 1.6271<br>1.4641 | 1.5700<br>1.4793 | 1.5350<br>1.4835 | 1.5516 | Ave | | 1.5311 | | | | 3.5 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48 Calibration End Date: 10/02/2015 14:04 Calibration ID: 30980

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2-Methylphenol | 1.4114 | 1.4902<br>1.3491 | 1.4012<br>1.3603 | 1.4223<br>1.3516 | 1.5104 | Ave | | 1.4121 | | | 0.7000 | 4.3 | | 20.0 | | | |
| bis (2-chloroisopropyl) ether | 2.1032 | 2.4010<br>1.9023 | 2.3063<br>1.8708 | 2.2908<br>1.8368 | 2.2583 | Ave | | 2.1212 | | | | 10.6 | | 20.0 | | | |
| Indene | 2.5417 | 2.8532<br>2.3321 | 2.6823<br>2.2562 | 2.7190<br>2.1708 | 2.7151 | Ave | | 2.5338 | | | | 9.9 | | 20.0 | | | |
| 3 & 4 Methylphenol | 1.4717 | 1.5716<br>1.3921 | 1.4945<br>1.3910 | 1.4916<br>1.3913 | 1.5495 | Ave | | 1.4692 | | | | 4.9 | | 20.0 | | | |
| N-Nitrosodi-n-propylamine | 0.9027 | 0.9251<br>0.8420 | 0.9104<br>0.8191 | 0.8990<br>0.7921 | 0.9376 | Ave | | 0.8785 | | | 0.5000 | 6.1 | | 20.0 | | | |
| Acetophenone | 2.0138 | 2.0766<br>1.8967 | 2.0491<br>1.8871 | 2.0752<br>1.8264 | 2.0787 | Ave | | 1.9880 | | | 0.0100 | 5.1 | | 20.0 | | | |
| Hexachloroethane | 0.6506 | 0.6511<br>0.6153 | 0.6147<br>0.6272 | 0.6386<br>0.6127 | 0.6770 | Ave | | 0.6359 | | | 0.3000 | 3.6 | | 20.0 | | | |
| Nitrobenzene | 0.3700 | 0.3890<br>0.3608 | 0.3813<br>0.3557 | 0.3686<br>0.3512 | 0.3786 | Ave | | 0.3694 | | | 0.2000 | 3.6 | | 20.0 | | | |
| Isophorone | 0.6883 | 0.7238<br>0.6711 | 0.6929<br>0.6633 | 0.6847<br>0.6524 | 0.7054 | Ave | | 0.6852 | | | 0.4000 | 3.4 | | 20.0 | | | |
| 2-Nitrophenol | 0.1876 | 0.1380<br>0.1801 | 0.1435<br>0.1805 | 0.1564<br>0.1754 | 0.1739 | Ave | | 0.1669 | | | 0.1000 | 11.1 | | 20.0 | | | |
| 2,4-Dimethylphenol | 0.3616 | 0.3894<br>0.3430 | 0.3809<br>0.3374 | 0.3682<br>0.3237 | 0.3597 | Ave | | 0.3580 | | | 0.2000 | 6.2 | | 20.0 | | | |
| Benzoic acid | 0.2291 | +++++<br>0.2141 | 0.0903<br>0.2178 | 0.1351<br>0.2228 | 0.1906 | Lin1 | -0.305 | 0.2283 | | | | | | | 0.9990 | | 0.9900 |
| Bis(2-chloroethoxy)methane | 0.4177 | 0.4647<br>0.3986 | 0.4508<br>0.3937 | 0.4375<br>0.3846 | 0.4412 | Ave | | 0.4236 | | | 0.3000 | 6.9 | | 20.0 | | | |
| 2,4-Dichlorophenol | 0.3004 | 0.3102<br>0.2890 | 0.2958<br>0.2884 | 0.2907<br>0.2810 | 0.2969 | Ave | | 0.2940 | | | 0.2000 | 3.0 | | 20.0 | | | |
| 1,2,4-Trichlorobenzene | 0.3054 | 0.3141<br>0.2927 | 0.3209<br>0.2858 | 0.3095<br>0.2778 | 0.3117 | Ave | | 0.3022 | | | | 5.0 | | 20.0 | | | |
| Naphthalene | 1.3004<br>1.0690 | 1.1308<br>1.0304 | 1.0735<br>1.0093 | 1.0684<br>0.9772 | 1.0969 | Ave | | 1.0840 | | | 0.7000 | 8.6 | | 20.0 | | | |
| 4-Chloroaniline | 0.4741 | 0.4930<br>0.4559 | 0.5069<br>0.4525 | 0.4937<br>0.4485 | 0.4800 | Ave | | 0.4756 | | | 0.0100 | 4.6 | | 20.0 | | | |
| 2,6-Dichlorophenol | 0.2958 | 0.3215<br>0.2856 | 0.3004<br>0.2833 | 0.2958<br>0.2781 | 0.2986 | Ave | | 0.2949 | | | | 4.5 | | 20.0 | | | |
| Hexachlorobutadiene | 0.1890 | 0.2046<br>0.1802 | 0.1919<br>0.1770 | 0.1878<br>0.1720 | 0.1923 | Ave | | 0.1868 | | | 0.0100 | 5.5 | | 20.0 | | | |
| Caprolactam | 0.1158 | 0.1186<br>0.1145 | 0.1187<br>0.1160 | 0.1123<br>0.1173 | 0.1180 | Ave | | 0.1164 | | | 0.0100 | 1.9 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48 Calibration End Date: 10/02/2015 14:04 Calibration ID: 30980

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 4-Chloro-3-methylphenol | 0.3081 | 0.3124<br>0.2963 | 0.3078<br>0.2958 | 0.3032<br>0.2895 | 0.3130 | Ave | | 0.3033 | | | 0.2000 | 2.8 | | 20.0 | | | |
| 2-Methylnaphthalene | 0.7756<br>0.6955 | 0.7490<br>0.6693 | 0.7170<br>0.6556 | 0.6957<br>0.6313 | 0.7165 | Ave | | 0.7006 | | | 0.4000 | 6.5 | | 20.0 | | | |
| 1-Methylnaphthalene | 0.7676<br>0.6107 | 0.6360<br>0.5848 | 0.6456<br>0.5768 | 0.6162<br>0.5626 | 0.6282 | Ave | | 0.6254 | | | | 9.6 | | 20.0 | | | |
| Hexachlorocyclopentadiene | 0.2948 | 0.2072<br>0.2918 | 0.2317<br>0.2918 | 0.2528<br>0.2873 | 0.2937 | Ave | | 0.2689 | | | 0.0500 | 12.7 | | 20.0 | | | |
| 1,2,4,5-Tetrachlorobenzene | 0.4885 | 0.5652<br>0.4629 | 0.5402<br>0.4463 | 0.5472<br>0.4297 | 0.5183 | Ave | | 0.4998 | | | 0.0100 | 10.1 | | 20.0 | | | |
| 2,4,6-Trichlorophenol | 0.3715 | 0.3724<br>0.3588 | 0.3515<br>0.3509 | 0.3712<br>0.3425 | 0.3759 | Ave | | 0.3619 | | | 0.2000 | 3.5 | | 20.0 | | | |
| 2,4,5-Trichlorophenol | 0.3894 | 0.3736<br>0.3765 | 0.3808<br>0.3754 | 0.3750<br>0.3713 | 0.3936 | Ave | | 0.3794 | | | 0.2000 | 2.1 | | 20.0 | | | |
| 1,1'-Biphenyl | 1.6748<br>1.3725 | 1.4785<br>1.3090 | 1.4844<br>1.2882 | 1.4329<br>1.2449 | 1.3992 | Ave | | 1.4094 | | | 0.0100 | 9.2 | | 20.0 | | | |
| 2-Chloronaphthalene | 1.4430<br>1.1165 | 1.3094<br>1.0904 | 1.1919<br>1.0799 | 1.2036<br>1.0476 | 1.1544 | Ave | | 1.1819 | | | 0.8000 | 10.7 | | 20.0 | | | |
| 2-Nitroaniline | 0.3659 | 0.3458<br>0.3517 | 0.3335<br>0.3522 | 0.3394<br>0.3426 | 0.3714 | Ave | | 0.3503 | | | 0.0100 | 3.7 | | 20.0 | | | |
| Dimethyl phthalate | 1.2918 | 1.3757<br>1.2433 | 1.3134<br>1.2373 | 1.2978<br>1.2027 | 1.3157 | Ave | | 1.2847 | | | 0.0100 | 4.3 | | 20.0 | | | |
| 1,3-Dinitrobenzene | 0.2128 | 0.1253<br>0.2137 | 0.1311<br>0.2169 | 0.1611<br>0.2086 | 0.1876 | Lin1 | -0.066 | 0.2149 | | | | | | | 0.9980 | | 0.9900 |
| 2,6-Dinitrotoluene | 0.2980 | 0.2441<br>0.2879 | 0.2649<br>0.2943 | 0.2864<br>0.2823 | 0.2962 | Ave | | 0.2818 | | | | 6.6 | | 20.0 | | | |
| Acenaphthylene | 2.0743<br>1.7077 | 1.8368<br>1.6563 | 1.7806<br>1.6705 | 1.7896<br>1.5999 | 1.7512 | Ave | | 1.7630 | | | 0.9000 | 7.8 | | 20.0 | | | |
| 3-Nitroaniline | 0.3380 | 0.2955<br>0.3324 | 0.3068<br>0.3328 | 0.3220<br>0.3280 | 0.3359 | Ave | | 0.3239 | | | 0.0100 | 4.7 | | 20.0 | | | |
| 2,4-Dinitrophenol | 0.1481 | +++++<br>0.1538 | 0.0351<br>0.1544 | 0.0510<br>0.1462 | 0.0918 | Lin1 | -0.328 | 0.1574 | | | 0.0100 | | | | 0.9920 | | 0.9900 |
| Acenaphthene | 1.3039<br>1.0225 | 1.2131<br>0.9782 | 1.1513<br>0.9499 | 1.1017<br>0.8846 | 1.0737 | Ave | | 1.0754 | | | 0.9000 | 12.4 | | 20.0 | | | |
| 4-Nitrophenol | 0.1794 | 0.1547<br>0.1768 | 0.1585<br>0.1774 | 0.1663<br>0.1698 | 0.1787 | Ave | | 0.1702 | | | | 5.6 | | 20.0 | | | |
| 2,4-Dinitrotoluene | 0.4036 | 0.2963<br>0.3894 | 0.3270<br>0.3928 | 0.3557<br>0.3847 | 0.3830 | Ave | | 0.3666 | | | 0.2000 | 10.2 | | 20.0 | | | |
| Dibenzofuran | 1.9069<br>1.5721 | 1.7236<br>1.5073 | 1.6855<br>1.5083 | 1.6327<br>1.4681 | 1.6051 | Ave | | 1.6233 | | | 0.8000 | 8.4 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48 Calibration End Date: 10/02/2015 14:04 Calibration ID: 30980

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2,3,4,6-Tetrachlorophenol | 0.3261 | 0.3032<br>0.3202 | 0.3101<br>0.3234 | 0.3209<br>0.3127 | 0.3268 | Ave | | 0.3179 | | | 0.0100 | 2.6 | | 20.0 | | | |
| Diethyl phthalate | 1.1755 | 1.3606<br>1.0860 | 1.2751<br>1.0391 | 1.2561<br>0.9672 | 1.2198 | Ave | | 1.1724 | | | 0.0100 | 11.3 | | 20.0 | | | |
| Hexadecane | 0.5384 | 0.6539<br>0.4892 | 0.6552<br>0.4399 | 0.6420<br>0.3964 | 0.6215 | Ave | | 0.5546 | | | | 18.6 | | 20.0 | | | |
| 4-Chlorophenyl phenyl ether | 0.6018 | 0.6868<br>0.5811 | 0.6476<br>0.5764 | 0.6229<br>0.5609 | 0.6109 | Ave | | 0.6110 | | | 0.4000 | 6.8 | | 20.0 | | | |
| 4-Nitroaniline | 0.3283 | 0.3168<br>0.3054 | 0.3309<br>0.2952 | 0.3259<br>0.2745 | 0.3421 | Ave | | 0.3149 | | | 0.0100 | 7.0 | | 20.0 | | | |
| Fluorene | 1.5331<br>1.3077 | 1.4393<br>1.2410 | 1.3917<br>1.1745 | 1.3621<br>1.1146 | 1.3380 | Ave | | 1.3224 | | | 0.9000 | 9.9 | | 20.0 | | | |
| 4,6-Dinitro-2-methylphenol | 0.1405 | +++++<br>0.1420 | 0.0535<br>0.1473 | 0.0736<br>0.1406 | 0.1077 | Lin1 | -0.238 | 0.1483 | | | 0.0100 | | | | 0.9970 | | 0.9900 |
| N-Nitrosodiphenylamine | 0.5721 | 0.6386<br>0.5413 | 0.6032<br>0.5217 | 0.6014<br>0.4871 | 0.6027 | Ave | | 0.5710 | | | 0.0100 | 8.9 | | 20.0 | | | |
| 1,2-Diphenylhydrazine(as Azobenzene) | 0.7979 | 0.8327<br>0.7726 | 0.8253<br>0.7603 | 0.8422<br>0.7222 | 0.8418 | Ave | | 0.7997 | | | | 5.5 | | 20.0 | | | |
| Azobenzene | 0.7979 | 0.8316<br>0.7731 | 0.8275<br>0.7602 | 0.8424<br>0.7228 | 0.8424 | Ave | | 0.7997 | | | | 5.5 | | 20.0 | | | |
| 4-Bromophenyl phenyl ether | 0.2291 | 0.2468<br>0.2220 | 0.2422<br>0.2182 | 0.2340<br>0.2104 | 0.2360 | Ave | | 0.2298 | | | 0.1000 | 5.4 | | 20.0 | | | |
| Hexachlorobenzene | 0.2897<br>0.2474 | 0.2906<br>0.2387 | 0.2639<br>0.2357 | 0.2702<br>0.2283 | 0.2592 | Ave | | 0.2582 | | | 0.1000 | 8.8 | | 20.0 | | | |
| Atrazine | 0.2085 | 0.2470<br>0.1989 | 0.2352<br>0.1914 | 0.2302<br>0.1784 | 0.2226 | Ave | | 0.2140 | | | 0.0100 | 11.1 | | 20.0 | | | |
| Pentachlorophenol | 0.1566 | 0.1203<br>0.1461 | 0.1373<br>0.1435 | 0.1462<br>0.1377 | 0.1580 | Ave | | 0.1432 | | | 0.0500 | 8.4 | | 20.0 | | | |
| n-Octadecane | 0.7811 | 1.0785<br>0.6664 | 1.0384<br>0.6036 | 1.0135<br>0.5462 | 0.9376 | Qua | 0.2920 | 0.8735 | -0.013823 | | | | | | 0.9980 | | 0.9900 |
| Phenanthrene | 1.3748<br>1.1479 | 1.2369<br>1.1131 | 1.1926<br>1.0919 | 1.1939<br>1.0530 | 1.1921 | Ave | | 1.1773 | | | 0.7000 | 8.0 | | 20.0 | | | |
| Anthracene | 1.3205<br>1.1782 | 1.2586<br>1.1423 | 1.2160<br>1.1412 | 1.1785<br>1.0647 | 1.2105 | Ave | | 1.1901 | | | 0.7000 | 6.2 | | 20.0 | | | |
| Carbazole | 1.1072 | 1.2431<br>1.0787 | 1.2009<br>1.0743 | 1.1353<br>1.0193 | 1.1600 | Ave | | 1.1273 | | | 0.0100 | 6.5 | | 20.0 | | | |
| Di-n-butyl phthalate | 1.3947 | 1.4647<br>1.3496 | 1.4241<br>1.3237 | 1.4029<br>1.2334 | 1.4453 | Ave | | 1.3798 | | | 0.0100 | 5.5 | | 20.0 | | | |
| Fluoranthene | 1.5321<br>1.2882 | 1.3939<br>1.2469 | 1.3537<br>1.2340 | 1.3252<br>1.1725 | 1.3424 | Ave | | 1.3210 | | | 0.6000 | 7.9 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48 Calibration End Date: 10/02/2015 14:04 Calibration ID: 30980

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Benzidine | 0.6451 | 0.5338<br>0.6050 | 0.5417<br>0.5945 | 0.5546<br>0.5565 | 0.6428 | Ave | | 0.5843 | | | | 7.6 | | 20.0 | | | |
| Pyrene | 1.3418<br>1.1061 | 1.1896<br>1.0765 | 1.1446<br>1.0713 | 1.1384<br>1.0412 | 1.1040 | Ave | | 1.1348 | | | 0.6000 | 7.9 | | 20.0 | | | |
| Butyl benzyl phthalate | 0.5176 | 0.5409<br>0.5064 | 0.5195<br>0.5022 | 0.5221<br>0.4846 | 0.5206 | Ave | | 0.5142 | | | 0.0100 | 3.2 | | 20.0 | | | |
| 3,3'-Dichlorobenzidine | 0.3938 | 0.4187<br>0.3806 | 0.4084<br>0.3666 | 0.4148<br>0.3572 | 0.4218 | Ave | | 0.3952 | | | 0.0100 | 6.3 | | 20.0 | | | |
| Bis(2-ethylhexyl) phthalate | 0.6563 | 0.7249<br>0.6056 | 0.7320<br>0.5577 | 0.7423<br>0.5175 | 0.7272 | Ave | | 0.6579 | | | 0.0100 | 13.4 | | 20.0 | | | |
| Benzo[a]anthracene | 1.2994<br>1.1072 | 1.1154<br>1.0880 | 1.1109<br>1.1166 | 1.1030<br>1.0814 | 1.1126 | Ave | | 1.1261 | | | 0.8000 | 5.9 | | 20.0 | | | |
| Chrysene | 1.2630<br>1.0330 | 1.1264<br>1.0316 | 1.0723<br>1.0239 | 1.0588<br>1.0141 | 1.0605 | Ave | | 1.0759 | | | 0.7000 | 7.2 | | 20.0 | | | |
| Di-n-octyl phthalate | 1.4250 | 1.2463<br>1.3943 | 1.3097<br>1.4235 | 1.3592<br>1.3602 | 1.4489 | Ave | | 1.3709 | | | 0.0100 | 4.9 | | 20.0 | | | |
| Benzo[b]fluoranthene | 1.1428<br>1.1817 | 1.1483<br>1.0957 | 1.1239<br>1.1339 | 1.0694<br>1.1422 | 1.1127 | Ave | | 1.1278 | | | 0.7000 | 2.9 | | 20.0 | | | |
| Benzo[k]fluoranthene | 1.4480<br>1.2152 | 1.3389<br>1.2296 | 1.3013<br>1.2500 | 1.2492<br>1.1684 | 1.2686 | Ave | | 1.2743 | | | 0.7000 | 6.4 | | 20.0 | | | |
| Benzo[a]pyrene | 1.1242<br>1.1335 | 1.1003<br>1.1221 | 1.1316<br>1.1430 | 1.0872<br>1.1046 | 1.1374 | Ave | | 1.1204 | | | 0.7000 | 1.7 | | 20.0 | | | |
| Indeno[1,2,3-cd]pyrene | 1.0594<br>1.2888 | 1.1682<br>1.2677 | 1.2083<br>1.2995 | 1.2340<br>1.2525 | 1.3208 | Ave | | 1.2332 | | | 0.5000 | 6.5 | | 20.0 | | | |
| Dibenz(a,h)anthracene | 0.7887<br>1.0727 | 1.0310<br>1.0681 | 1.0191<br>1.0966 | 1.0137<br>1.0546 | 1.0748 | Ave | | 1.0244 | | | 0.4000 | 9.0 | | 20.0 | | | |
| Benzo[g,h,i]perylene | 0.9436<br>1.1332 | 1.0999<br>1.1069 | 1.0654<br>1.1485 | 1.0731<br>1.1229 | 1.1388 | Ave | | 1.0925 | | | 0.5000 | 5.7 | | 20.0 | | | |
| 2-Fluorophenol (Surr) | 1.4936 | 1.4868<br>1.4489 | 1.4884<br>1.4663 | 1.5162<br>1.4709 | 1.5176 | Ave | | 1.4861 | | | | 1.6 | | 20.0 | | | |
| Phenol-d5 (Surr) | 1.8356 | 1.9501<br>1.7490 | 1.8116<br>1.7382 | 1.7830<br>1.7442 | 1.8996 | Ave | | 1.8139 | | | | 4.3 | | 20.0 | | | |
| Nitrobenzene-d5 (Surr) | 0.4177<br>0.3995 | 0.4058<br>0.3874 | 0.3992<br>0.3905 | 0.3968<br>0.3818 | 0.4054 | Ave | | 0.3982 | | | | 2.7 | | 20.0 | | | |
| 2-Fluorobiphenyl (Surr) | 1.6344<br>1.2399 | 1.4109<br>1.1877 | 1.3265<br>1.1814 | 1.2886<br>1.1377 | 1.2630 | Ave | | 1.2967 | | | | 11.7 | | 20.0 | | | |
| 2,4,6-Tribromophenol (Surr) | 0.1986 | 0.1782<br>0.1920 | 0.1845<br>0.1971 | 0.1810<br>0.1944 | 0.1873 | Ave | | 0.1891 | | | | 4.0 | | 20.0 | | | |
| Terphenyl-d14 (Surr) | 0.9059<br>0.7324 | 0.8249<br>0.7190 | 0.7668<br>0.7171 | 0.7574<br>0.6951 | 0.7388 | Ave | | 0.7619 | | | | 8.6 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48 Calibration End Date: 10/02/2015 14:04 Calibration ID: 30980

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-200091/6 | 51002006.D |
| Level 2 | STD2 240-200091/5 | 51002005.D |
| Level 3 | STD3 240-200091/4 | 51002004.D |
| Level 4 | STD4 240-200091/3 | 51002003.D |
| Level 5 | STD5 240-200091/2 | 51002002.D |
| Level 6 | STD6 240-200091/10 | 51002010.D |
| Level 7 | STD7 240-200091/9 | 51002009.D |
| Level 8 | STD8 240-200091/8 | 51002008.D |
| Level 9 | STD9 240-200091/7 | 51002007.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1,4-Dioxane | DCB | Ave | 370662 | 20911<br>545261 | 37914<br>731845 | 68248<br>915986 | 166746 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| N-Nitrosodimethylamine | DCB | Ave | 510065 | 26906<br>763855 | 48095<br>1005940 | 91510<br>1260405 | 219541 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Pyridine | DCB | Ave | 927473 | 47995<br>1384286 | 86556<br>1858888 | 164774<br>2312079 | 409338 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzaldehyde | DCB | Ave | 1235862 | 68753<br>1680758 | 130916<br>2138376 | 244675<br>2426096 | 566513 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Phenol | DCB | Ave | 1039342 | 55897<br>1520706 | 101626<br>2004194 | 195041<br>2436933 | 455561 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Aniline | DCB | Ave | 1343901 | 67917<br>1953885 | 128107<br>2583026 | 241843<br>3204478 | 578515 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Bis(2-chloroethyl)ether | DCB | Ave | 843878 | 43637<br>1220540 | 81176<br>1603524 | 152518<br>1945416 | 364177 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Chlorophenol | DCB | Ave | 854637 | 43227<br>1262124 | 84781<br>1676912 | 156548<br>2033603 | 365089 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| n-Decane | DCB | Ave | 1041214 | 57768<br>1506688 | 109265<br>1941972 | 197615<br>2328538 | 472437 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,3-Dichlorobenzene | DCB | Ave | 881944 | 45808<br>1268801 | 83623<br>1674644 | 163103<br>2039987 | 385698 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,4-Dichlorobenzene | DCB | Ave | 901675 | 47174<br>1291211 | 87875<br>1726579 | 161346<br>2081257 | 393078 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzyl alcohol | DCB | Ave | 567465 | 30245<br>827651 | 54243<br>1101205 | 102751<br>1338352 | 248213 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2-Dichlorobenzene | DCB | Ave | 851645 | 44397<br>1237679 | 83698<br>1640215 | 154250<br>1999182 | 363748 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Methylphenol | DCB | Ave | 781451 | 40662<br>1140453 | 74700<br>1508237 | 142925<br>1821326 | 354108 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48 Calibration End Date: 10/02/2015 14:04 Calibration ID: 30980

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| bis (2-chloroisopropyl) ether | DCB | Ave | 1164480 | 65511<br>1608103 | 122954<br>2074273 | 230203<br>2475238 | 529431 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Indene | DCB | Ave | 2814476 | 155700<br>3942904 | 285997<br>5003192 | 546467<br>5850511 | 1273077 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 3 & 4 Methylphenol | DCB | Ave | 814794 | 42882<br>1176808 | 79676<br>1542266 | 149893<br>1874933 | 363254 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| N-Nitrosodi-n-propylamine | DCB | Ave | 499813 | 25243<br>711760 | 48533<br>908131 | 90336<br>1067435 | 219816 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Acetophenone | DCB | Ave | 1114952 | 56662<br>1603405 | 109240<br>2092353 | 208536<br>2461240 | 487336 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachloroethane | DCB | Ave | 360194 | 17766<br>520125 | 32770<br>695444 | 64178<br>825687 | 158706 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Nitrobenzene | NPT | Ave | 838541 | 43092<br>1234105 | 80768<br>1631735 | 151101<br>1987886 | 364359 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Isophorone | NPT | Ave | 1559881 | 80173<br>2295589 | 146792<br>3043071 | 280704<br>3692856 | 678797 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Nitrophenol | NPT | Ave | 425227 | 15291<br>615997 | 30400<br>828256 | 64101<br>992550 | 167320 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dimethylphenol | NPT | Ave | 819350 | 43130<br>1173285 | 80689<br>1547760 | 150936<br>1832271 | 346175 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzoic acid | NPT | Lin1 | 1038142 | +++++<br>1464610 | 38247<br>1998568 | 110811<br>2522170 | 366851 | 20.0 | +++++<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Bis(2-chloroethoxy)methane | NPT | Ave | 946659 | 51471<br>1363574 | 95507<br>1805907 | 179349<br>2176975 | 424597 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dichlorophenol | NPT | Ave | 680679 | 34360<br>988632 | 62666<br>1322854 | 119194<br>1590608 | 285685 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2,4-Trichlorobenzene | NPT | Ave | 692195 | 34798<br>1001246 | 67981<br>1310983 | 126903<br>1572169 | 299921 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Naphthalene | NPT | Ave | 28740<br>2422576 | 125257<br>3524618 | 227409<br>4629906 | 438031<br>5531113 | 1055573 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Chloroaniline | NPT | Ave | 1074448 | 54612<br>1559338 | 107392<br>2075944 | 202393<br>2538434 | 461884 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,6-Dichlorophenol | NPT | Ave | 670420 | 35610<br>976864 | 63644<br>1299712 | 121289<br>1574160 | 287306 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorobutadiene | NPT | Ave | 428273 | 22658<br>616384 | 40647<br>811900 | 76997<br>973311 | 185013 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Caprolactam | NPT | Ave | 524922 | 26271<br>783569 | 50274<br>1064258 | 92048<br>1328164 | 227125 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 4-Chloro-3-methylphenol | NPT | Ave | 698207 | 34610<br>1013626 | 65210<br>1356946 | 124311<br>1638397 | 301160 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Methylnaphthalene | NPT | Ave | 17141<br>1576062 | 82967<br>2289268 | 151895<br>3007482 | 285213<br>3573068 | 689536 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48 Calibration End Date: 10/02/2015 14:04 Calibration ID: 30980

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1-Methylnaphthalene | NPT | Ave | 16966<br>1383973 | 70455<br>2000288 | 136756<br>2645906 | 252647<br>3184434 | 604477 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorocyclopentadiene | ANT | Ave | <br>390085 | 12564<br>584683 | 27556<br>777885 | 57197<br>948047 | 161161 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2,4,5-Tetrachlorobenzene | ANT | Ave | <br>646406 | 34273<br>927687 | 64237<br>1189970 | 123837<br>1418071 | 284441 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,6-Trichlorophenol | ANT | Ave | <br>491572 | 22581<br>719102 | 41801<br>935674 | 83990<br>1130369 | 206296 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,5-Trichlorophenol | ANT | Ave | <br>515195 | 22656<br>754448 | 45283<br>1000757 | 84859<br>1225297 | 216019 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,1'-Biphenyl | ANT | Ave | 20494<br>1816044 | 89656<br>2623309 | 176516<br>3434479 | 324251<br>4108549 | 767845 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Chloronaphthalene | ANT | Ave | 17658<br>1477280 | 79399<br>2185216 | 141725<br>2879241 | 272377<br>3457566 | 633504 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Nitroaniline | ANT | Ave | <br>484150 | 20971<br>704890 | 39662<br>938975 | 76808<br>1130855 | 203833 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dimethyl phthalate | ANT | Ave | <br>1709256 | 83422<br>2491597 | 156176<br>3298887 | 293685<br>3969492 | 722024 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,3-Dinitrobenzene | ANT | Lin1 | <br>281574 | 7597<br>428331 | 15586<br>578326 | 36454<br>688473 | 102927 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,6-Dinitrotoluene | ANT | Ave | <br>394339 | 14804<br>576932 | 31503<br>784639 | 64800<br>931549 | 162540 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Acenaphthylene | ANT | Ave | 25382<br>2259522 | 111380<br>3319231 | 211730<br>4453755 | 404966<br>5280391 | 960986 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 3-Nitroaniline | ANT | Ave | <br>447285 | 17920<br>666109 | 36480<br>887185 | 72876<br>1082650 | 184318 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dinitrophenol | ANT | Lin1 | <br>391802 | +++++<br>616389 | 8344<br>823125 | 23069<br>964987 | 100738 | <br>20.0 | +++++<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Acenaphthene | ANT | Ave | 15956<br>1352877 | 73560<br>1960267 | 136902<br>2532651 | 249312<br>2919429 | 589216 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Nitrophenol | ANT | Ave | <br>474712 | 18758<br>708462 | 37701<br>945992 | 75258<br>1120709 | 196147 | <br>20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 2,4-Dinitrotoluene | ANT | Ave | <br>534050 | 17970<br>780449 | 38883<br>1047168 | 80488<br>1269724 | 210162 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dibenzofuran | ANT | Ave | 23334<br>2080160 | 104518<br>3020592 | 200424<br>4021214 | 369474<br>4845200 | 880799 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,3,4,6-Tetrachlorophenol | ANT | Ave | <br>431493 | 18383<br>641722 | 36874<br>862275 | 72619<br>1032006 | 179341 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Diethyl phthalate | ANT | Ave | <br>1555322 | 82503<br>2176407 | 151628<br>2770424 | 284239<br>3191986 | 669397 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexadecane | PHN | Ave | <br>1131032 | 62955<br>1546902 | 123571<br>1870574 | 231460<br>2114114 | 530812 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48 Calibration End Date: 10/02/2015 14:04 Calibration ID: 30980

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 4-Chlorophenyl phenyl ether | ANT | Ave | 796301 | 41646<br>1164488 | 77006<br>1536749 | 140959<br>1851206 | 335228 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Nitroaniline | ANT | Ave | 434429 | 19209<br>611969 | 39345<br>787092 | 73747<br>905845 | 187751 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Fluorene | ANT | Ave | 18760<br>1730223 | 87274<br>2486967 | 165484<br>3131335 | 308240<br>3678529 | 734219 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4,6-Dinitro-2-methylphenol | PHN | Lin1 | 590204 | +++++<br>897806 | 20166<br>1253065 | 53043<br>1499930 | 183900 | 20.0 | +++++<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| N-Nitrosodiphenylamine | PHN | Ave | 2403332 | 122960<br>3423294 | 227499<br>4437465 | 433673<br>5196185 | 1029508 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | PHN | Ave | 1675941 | 80170<br>2443006 | 155647<br>3233352 | 303647<br>3852013 | 718987 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Azobenzene | PHN | Ave | 1676087 | 80066<br>2444390 | 156051<br>3232944 | 303700<br>3855135 | 719496 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Bromophenyl phenyl ether | PHN | Ave | 481209 | 23757<br>701931 | 45684<br>928120 | 84352<br>1122446 | 201544 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorobenzene | PHN | Ave | 5609<br>519703 | 27980<br>754663 | 49774<br>1002347 | 97403<br>1217956 | 221331 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Atrazine | PHN | Ave | 875985 | 47563<br>1257802 | 88712<br>1627481 | 165998<br>1902732 | 380168 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Pentachlorophenol | PHN | Ave | 657894 | 23164<br>924064 | 51777<br>1220147 | 105393<br>1469132 | 269822 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| n-Octadecane | ANT | Qua | 1033464 | 65400<br>1335536 | 123473<br>1609282 | 229344<br>1802677 | 514499 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Phenanthrene | PHN | Ave | 26616<br>2411126 | 119089<br>3519639 | 224901<br>4643298 | 430431<br>5616503 | 1018122 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Anthracene | PHN | Ave | 25564<br>2474768 | 121176<br>3611735 | 229322<br>4853200 | 424899<br>5679119 | 1033829 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Carbazole | PHN | Ave | 2325742 | 119681<br>3410796 | 226468<br>4568414 | 409309<br>5436592 | 990732 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Di-n-butyl phthalate | PHN | Ave | 2929632 | 141017<br>4267270 | 268573<br>5629384 | 505808<br>6579074 | 1234410 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Fluoranthene | PHN | Ave | 29660<br>2705932 | 134205<br>3942504 | 255297<br>5247887 | 477790<br>6253971 | 1146489 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzidine | CRY | Ave | 3345321 | 126524<br>4745755 | 251450<br>6239797 | 488303<br>7217928 | 1372140 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Pyrene | CRY | Ave | 31197<br>2867797 | 140967<br>4222040 | 265647<br>5621799 | 501177<br>6752901 | 1178311 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Butyl benzyl phthalate | CRY | Ave | 1342117 | 64093<br>1986008 | 120560<br>2635406 | 229845<br>3143034 | 555638 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 3,3'-Dichlorobenzidine | CRY | Ave | 2042036 | 99232<br>2985002 | 189560<br>3847725 | 365246<br>4633511 | 900335 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48 Calibration End Date: 10/02/2015 14:04 Calibration ID: 30980

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Bis(2-ethylhexyl) phthalate | CRY | Ave | 1701558 | 85896<br>2374966 | 169877<br>2926383 | 326796<br>3356175 | 776158 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[a]anthracene | CRY | Ave | 30212<br>2870719 | 132181<br>4266960 | 257826<br>5859665 | 485590<br>7013789 | 1187464 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Chrysene | CRY | Ave | 29365<br>2678230 | 133481<br>4045667 | 248867<br>5373291 | 466114<br>6576826 | 1131795 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Di-n-octyl phthalate | PRY | Ave | 3402387 | 133560<br>5122374 | 277640<br>6897735 | 556183<br>8332711 | 1431617 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[b]fluoranthene | PRY | Ave | 24087<br>2821318 | 123058<br>4025307 | 238255<br>5494463 | 437589<br>6997425 | 1099354 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[k]fluoranthene | PRY | Ave | 30521<br>2901469 | 143482<br>4517162 | 275861<br>6056952 | 511155<br>7157600 | 1253398 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[a]pyrene | PRY | Ave | 23695<br>2706410 | 117909<br>4122371 | 239888<br>5538764 | 444893<br>6767133 | 1123767 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Indeno[1,2,3-cd]pyrene | PRY | Ave | 22329<br>3077152 | 125185<br>4657108 | 256153<br>6296984 | 504951<br>7672812 | 1304973 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dibenz(a,h)anthracene | PRY | Ave | 16624<br>2561245 | 110491<br>3923783 | 216043<br>5313779 | 414824<br>6460507 | 1061935 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[g,h,i]perylene | PRY | Ave | 19890<br>2705601 | 117865<br>4066451 | 225855<br>5565214 | 439115<br>6878912 | 1125146 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Fluorophenol (Surr) | DCB | Ave | 826968 | 40568<br>1224794 | 79350<br>1625797 | 152367<br>1982168 | 355783 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Phenol-d5 (Surr) | DCB | Ave | 1016291 | 53208<br>1478509 | 96582<br>1927273 | 179175<br>2350446 | 445346 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Nitrobenzene-d5 (Surr) | NPT | Ave | 9232<br>905450 | 44953<br>1324964 | 84570<br>1791496 | 162679<br>2161035 | 390090 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Fluorobiphenyl (Surr) | ANT | Ave | 20000<br>1640609 | 85555<br>2380059 | 157732<br>3149652 | 291601<br>3754930 | 693115 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,6-Tribromophenol (Surr) | ANT | Ave | 262754 | 10804<br>384804 | 21934<br>525506 | 40962<br>641586 | 102778 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Terphenyl-d14 (Surr) | CRY | Ave | 21063<br>1898847 | 97756<br>2819906 | 177962<br>3763219 | 333450<br>4507965 | 788542 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin1 = Linear 1/conc ISTD |
| Qua = Quadratic ISTD |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19  Calibration End Date: 10/30/2015 16:42  Calibration ID: 31170

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-204569/6 | 51030006.D |
| Level 2 | STD2 240-204569/5 | 51030005.D |
| Level 3 | STD3 240-204569/4 | 51030004.D |
| Level 4 | STD4 240-204569/3 | 51030003.D |
| Level 5 | STD5 240-204569/2 | 51030002.D |
| Level 6 | STD6 240-204569/10 | 51030010.D |
| Level 7 | STD7 240-204569/9 | 51030009.D |
| Level 8 | STD8 240-204569/8 | 51030008.D |
| Level 9 | STD9 240-204569/7 | 51030007.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,4-Dioxane | 0.7879 | 0.6702<br>0.7724 | 0.7963<br>0.7912 | 0.7519<br>0.7577 | 0.7576 | Ave | | 0.7606 | | | | 5.3 | | 20.0 | | | |
| N-Nitrosodimethylamine | 1.1590 | 1.0469<br>1.1257 | 0.9916<br>1.1673 | 1.0993<br>1.1059 | 1.0809 | Ave | | 1.0971 | | | | 5.3 | | 20.0 | | | |
| Pyridine | 2.0907 | 2.0205<br>1.9910 | 1.8876<br>2.0977 | 2.0118<br>1.9855 | 2.0565 | Ave | | 2.0177 | | | | 3.3 | | 20.0 | | | |
| Benzaldehyde | 1.3725 | 1.5727<br>1.2076 | 1.5657<br>1.2344 | 1.5133<br>+++++ | 1.4739 | Ave | | 1.4200 | | | 0.0100 | 10.7 | | 20.0 | | | |
| Phenol | 2.1853 | 2.1727<br>2.0136 | 2.0934<br>2.0804 | 2.1195<br>1.9722 | 2.1525 | Ave | | 2.0987 | | | 0.8000 | 3.6 | | 20.0 | | | |
| Aniline | 2.7280 | 2.6247<br>2.5385 | 2.6136<br>2.6466 | 2.6262<br>2.5459 | 2.6111 | Ave | | 2.6168 | | | | 2.3 | | 20.0 | | | |
| Bis(2-chloroethyl)ether | 1.8558 | 1.9142<br>1.6942 | 1.8429<br>1.7572 | 1.8447<br>1.6890 | 1.8270 | Ave | | 1.8031 | | | 0.7000 | 4.5 | | 20.0 | | | |
| 2-Chlorophenol | 1.5989 | 1.4425<br>1.5006 | 1.5455<br>1.5432 | 1.5087<br>1.4819 | 1.5120 | Ave | | 1.5167 | | | 0.8000 | 3.1 | | 20.0 | | | |
| n-Decane | 1.7829 | 1.8718<br>1.6389 | 1.8936<br>1.6552 | 1.8312<br>1.5772 | 1.8099 | Ave | | 1.7576 | | | | 6.7 | | 20.0 | | | |
| 1,3-Dichlorobenzene | 1.6084 | 1.5596<br>1.5011 | 1.6679<br>1.5578 | 1.6175<br>1.4910 | 1.5792 | Ave | | 1.5728 | | | | 3.8 | | 20.0 | | | |
| 1,4-Dichlorobenzene | 1.6334 | 1.6944<br>1.5231 | 1.6061<br>1.5860 | 1.6427<br>1.5227 | 1.6060 | Ave | | 1.6018 | | | | 3.7 | | 20.0 | | | |
| Benzyl alcohol | 1.1163 | 0.8791<br>1.0801 | 0.9185<br>1.1234 | 0.9897<br>1.0846 | 1.0413 | Ave | | 1.0291 | | | | 8.9 | | 20.0 | | | |
| 1,2-Dichlorobenzene | 1.5481 | 1.6647<br>1.4507 | 1.5548<br>1.5234 | 1.5786<br>1.4511 | 1.5515 | Ave | | 1.5404 | | | | 4.5 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2-Methylphenol | 1.6435 | 1.5529<br>1.5278 | 1.5884<br>1.6065 | 1.5262<br>1.5342 | 1.5781 | Ave | | 1.5697 | | | 0.7000 | 2.7 | | 20.0 | | | |
| bis (2-chloroisopropyl) ether | 2.0663 | 2.1290<br>1.9114 | 2.1504<br>1.9756 | 2.0501<br>1.8678 | 2.0388 | Ave | | 2.0237 | | | | 4.9 | | 20.0 | | | |
| Indene | 2.7276 | 2.8401<br>2.4658 | 2.8969<br>2.5378 | 2.8654<br>2.3993 | 2.7687 | Ave | | 2.6877 | | | | 7.2 | | 20.0 | | | |
| 3 & 4 Methylphenol | 1.6857 | 1.6121<br>1.5969 | 1.5331<br>1.6527 | 1.6625<br>1.5643 | 1.6490 | Ave | | 1.6195 | | | | 3.2 | | 20.0 | | | |
| N-Nitrosodi-n-propylamine | 1.1989 | 1.1171<br>1.1083 | 1.1602<br>1.1430 | 1.1470<br>1.0752 | 1.1583 | Ave | | 1.1385 | | | 0.5000 | 3.3 | | 20.0 | | | |
| Acetophenone | 2.3832 | 2.4667<br>2.2595 | 2.3410<br>2.3395 | 2.3767<br>2.2604 | 2.3981 | Ave | | 2.3531 | | | 0.0100 | 3.0 | | 20.0 | | | |
| Hexachloroethane | 0.6860 | 0.6898<br>0.6531 | 0.6582<br>0.6793 | 0.6592<br>0.6596 | 0.6614 | Ave | | 0.6683 | | | 0.3000 | 2.1 | | 20.0 | | | |
| Nitrobenzene | 0.4587 | 0.4485<br>0.4327 | 0.4414<br>0.4340 | 0.4609<br>0.4248 | 0.4662 | Ave | | 0.4459 | | | 0.2000 | 3.4 | | 20.0 | | | |
| Isophorone | 0.7828 | 0.7464<br>0.7590 | 0.7806<br>0.7476 | 0.7726<br>0.7157 | 0.7963 | Ave | | 0.7626 | | | 0.4000 | 3.4 | | 20.0 | | | |
| 2-Nitrophenol | 0.1828 | 0.1373<br>0.1790 | 0.1472<br>0.1753 | 0.1582<br>0.1700 | 0.1781 | Ave | | 0.1660 | | | 0.1000 | 10.0 | | 20.0 | | | |
| 2,4-Dimethylphenol | 0.3869 | 0.3987<br>0.3723 | 0.4012<br>0.3647 | 0.3813<br>0.3471 | 0.4020 | Ave | | 0.3818 | | | 0.2000 | 5.1 | | 20.0 | | | |
| Benzoic acid | 0.2127 | +++++<br>0.2174 | 0.0852<br>0.2213 | 0.1129<br>0.2114 | 0.1510 | Lin1 | -0.359 | 0.2234 | | | | | | | 0.9960 | | 0.9900 |
| Bis(2-chloroethoxy)methane | 0.4632 | 0.4830<br>0.4267 | 0.4845<br>0.4231 | 0.4695<br>0.4030 | 0.4779 | Ave | | 0.4539 | | | 0.3000 | 7.0 | | 20.0 | | | |
| 2,4-Dichlorophenol | 0.2586 | 0.2322<br>0.2510 | 0.2551<br>0.2491 | 0.2461<br>0.2424 | 0.2571 | Ave | | 0.2489 | | | 0.2000 | 3.5 | | 20.0 | | | |
| 1,2,4-Trichlorobenzene | 0.2904 | 0.3057<br>0.2778 | 0.3084<br>0.2753 | 0.2978<br>0.2688 | 0.3048 | Ave | | 0.2911 | | | | 5.3 | | 20.0 | | | |
| Naphthalene | 1.3148<br>1.0745 | 1.1643<br>1.0146 | 1.1361<br>0.9888 | 1.0999<br>0.9493 | 1.0983 | Ave | | 1.0934 | | | 0.7000 | 9.9 | | 20.0 | | | |
| 4-Chloroaniline | 0.4907 | 0.4694<br>0.4709 | 0.4675<br>0.4648 | 0.4499<br>0.4471 | 0.4606 | Ave | | 0.4651 | | | 0.0100 | 2.9 | | 20.0 | | | |
| 2,6-Dichlorophenol | 0.2619 | 0.2609<br>0.2504 | 0.2649<br>0.2440 | 0.2474<br>0.2401 | 0.2580 | Ave | | 0.2535 | | | | 3.6 | | 20.0 | | | |
| Hexachlorobutadiene | 0.1543 | 0.1757<br>0.1478 | 0.1596<br>0.1459 | 0.1568<br>0.1417 | 0.1563 | Ave | | 0.1548 | | | 0.0100 | 6.8 | | 20.0 | | | |
| Caprolactam | 0.1199 | 0.0914<br>0.1182 | 0.1059<br>0.1173 | 0.1129<br>0.1147 | 0.1141 | Ave | | 0.1118 | | | 0.0100 | 8.3 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 4-Chloro-3-methylphenol | 0.3405 | 0.2915<br>0.3291 | 0.2982<br>0.3289 | 0.3051<br>0.3200 | 0.3280 | Ave | | 0.3177 | | | 0.2000 | 5.5 | | 20.0 | | | |
| 2-Methylnaphthalene | 0.6714<br>0.6849 | 0.7552<br>0.6463 | 0.7151<br>0.6273 | 0.6901<br>0.5996 | 0.7011 | Ave | | 0.6768 | | | 0.4000 | 7.0 | | 20.0 | | | |
| 1-Methylnaphthalene | 0.6508<br>0.6069 | 0.6257<br>0.5675 | 0.6232<br>0.5547 | 0.6191<br>0.5412 | 0.6146 | Ave | | 0.6004 | | | | 6.2 | | 20.0 | | | |
| Hexachlorocyclopentadiene | 0.2836 | 0.2033<br>0.2937 | 0.2109<br>0.2923 | 0.2357<br>0.2863 | 0.2621 | Ave | | 0.2585 | | | 0.0500 | 14.3 | | 20.0 | | | |
| 1,2,4,5-Tetrachlorobenzene | 0.4829 | 0.5543<br>0.4681 | 0.5241<br>0.4467 | 0.5277<br>0.4361 | 0.5085 | Ave | | 0.4935 | | | 0.0100 | 8.5 | | 20.0 | | | |
| 2,4,6-Trichlorophenol | 0.3520 | 0.3202<br>0.3511 | 0.3243<br>0.3405 | 0.3380<br>0.3354 | 0.3397 | Ave | | 0.3377 | | | 0.2000 | 3.3 | | 20.0 | | | |
| 2,4,5-Trichlorophenol | 0.3611 | 0.3258<br>0.3643 | 0.3548<br>0.3514 | 0.3345<br>0.3507 | 0.3665 | Ave | | 0.3511 | | | 0.2000 | 4.1 | | 20.0 | | | |
| 1,1'-Biphenyl | 1.7382<br>1.5054 | 1.7110<br>1.4492 | 1.5822<br>1.3989 | 1.5786<br>1.3700 | 1.5456 | Ave | | 1.5421 | | | 0.0100 | 8.3 | | 20.0 | | | |
| 2-Chloronaphthalene | 1.3119<br>1.1298 | 1.2573<br>1.0949 | 1.1833<br>1.0732 | 1.1920<br>1.0363 | 1.1795 | Ave | | 1.1620 | | | 0.8000 | 7.6 | | 20.0 | | | |
| 2-Nitroaniline | 0.4424 | 0.3228<br>0.4442 | 0.3489<br>0.4281 | 0.3983<br>0.4198 | 0.4199 | Ave | | 0.4031 | | | 0.0100 | 11.0 | | 20.0 | | | |
| Dimethyl phthalate | 1.2250 | 1.3046<br>1.1901 | 1.2480<br>1.1306 | 1.2490<br>1.1063 | 1.2488 | Ave | | 1.2128 | | | 0.0100 | 5.5 | | 20.0 | | | |
| 1,3-Dinitrobenzene | 0.2148 | 0.1418<br>0.2208 | 0.1424<br>0.2112 | 0.1739<br>0.2124 | 0.1993 | Ave | | 0.1896 | | | | 17.2 | | 20.0 | | | |
| 2,6-Dinitrotoluene | 0.3015 | 0.2515<br>0.2974 | 0.2744<br>0.2833 | 0.2734<br>0.2819 | 0.2878 | Ave | | 0.2814 | | | | 5.6 | | 20.0 | | | |
| Acenaphthylene | 2.1239<br>1.9191 | 1.8953<br>1.8490 | 1.9145<br>1.7759 | 1.9143<br>1.7344 | 1.9293 | Ave | | 1.8951 | | | 0.9000 | 5.8 | | 20.0 | | | |
| 3-Nitroaniline | 0.3684 | 0.2911<br>0.3658 | 0.3323<br>0.3564 | 0.3421<br>0.3575 | 0.3522 | Ave | | 0.3457 | | | 0.0100 | 7.2 | | 20.0 | | | |
| 2,4-Dinitrophenol | 0.1226 | +++++<br>0.1490 | 0.0184<br>0.1495 | 0.0295<br>0.1498 | 0.0689 | Lin | -0.576 | 0.1620 | | | 0.0100 | | | | 0.9940 | | 0.9900 |
| Acenaphthene | 1.4818<br>1.1919 | 1.3172<br>1.1478 | 1.2378<br>1.0901 | 1.2469<br>1.0790 | 1.2360 | Ave | | 1.2254 | | | 0.9000 | 10.1 | | 20.0 | | | |
| 4-Nitrophenol | 0.2363 | 0.1350<br>0.2358 | 0.1725<br>0.2308 | 0.1994<br>0.2253 | 0.2085 | Ave | | 0.2054 | | | | 17.5 | | 20.0 | | | |
| 2,4-Dinitrotoluene | 0.3819 | 0.2653<br>0.3890 | 0.3195<br>0.3921 | 0.3356<br>0.3853 | 0.3775 | Ave | | 0.3558 | | | 0.2000 | 12.7 | | 20.0 | | | |
| Dibenzofuran | 1.9633<br>1.5887 | 1.8065<br>1.5423 | 1.6844<br>1.4828 | 1.6535<br>1.4457 | 1.6428 | Ave | | 1.6455 | | | 0.8000 | 9.8 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2,3,4,6-Tetrachlorophenol | 0.2882 | 0.2083<br>0.2752 | 0.2408<br>0.2694 | 0.2697<br>0.2698 | 0.2803 | Ave | | 0.2627 | | | 0.0100 | 9.9 | | 20.0 | | | |
| Hexadecane | 0.8423 | 0.9528<br>0.7419 | 0.9952<br>0.6898 | 0.9799<br>+++++ | 0.9285 | Ave | | 0.8758 | | | | 13.8 | | 20.0 | | | |
| Diethyl phthalate | 1.1435 | 1.3219<br>1.1140 | 1.3222<br>1.0714 | 1.3321<br>1.0617 | 1.2828 | Ave | | 1.2062 | | | 0.0100 | 9.9 | | 20.0 | | | |
| 4-Chlorophenyl phenyl ether | 0.5893 | 0.6305<br>0.5829 | 0.6266<br>0.5548 | 0.6099<br>0.5380 | 0.5949 | Ave | | 0.5909 | | | 0.4000 | 5.5 | | 20.0 | | | |
| 4-Nitroaniline | 0.3862 | 0.2821<br>0.3889 | 0.3158<br>0.3676 | 0.3459<br>0.3461 | 0.3410 | Ave | | 0.3467 | | | 0.0100 | 10.3 | | 20.0 | | | |
| Fluorene | 1.4143<br>1.3000 | 1.3977<br>1.2358 | 1.3185<br>1.1690 | 1.3302<br>1.1121 | 1.3108 | Ave | | 1.2876 | | | 0.9000 | 7.7 | | 20.0 | | | |
| 4,6-Dinitro-2-methylphenol | 0.1273 | +++++<br>0.1312 | 0.0456<br>0.1282 | 0.0664<br>0.1258 | 0.0946 | Lin1 | -0.215 | 0.1327 | | | 0.0100 | | | | 0.9970 | | 0.9900 |
| N-Nitrosodiphenylamine | 0.5686 | 0.5541<br>0.4887 | 0.5453<br>0.4764 | 0.5445<br>0.4597 | 0.5726 | Ave | | 0.5262 | | | 0.0100 | 8.4 | | 20.0 | | | |
| 1,2-Diphenylhydrazine(as Azobenzene) | 1.1353 | 1.0022<br>0.9923 | 0.9840<br>0.9576 | 1.0232<br>0.9032 | 1.1113 | Ave | | 1.0141 | | | | 7.6 | | 20.0 | | | |
| Azobenzene | 1.1352 | 1.0064<br>0.9924 | 0.9829<br>0.9577 | 1.0228<br>0.9034 | 1.1118 | Ave | | 1.0141 | | | | 7.6 | | 20.0 | | | |
| 4-Bromophenyl phenyl ether | 0.2076 | 0.1841<br>0.1792 | 0.1926<br>0.1792 | 0.1849<br>0.1705 | 0.1914 | Ave | | 0.1862 | | | 0.1000 | 6.0 | | 20.0 | | | |
| Hexachlorobenzene | 0.2326<br>0.2013 | 0.2025<br>0.1742 | 0.1926<br>0.1716 | 0.1883<br>0.1642 | 0.1879 | Ave | | 0.1906 | | | 0.1000 | 10.8 | | 20.0 | | | |
| Atrazine | 0.2016 | 0.1731<br>0.1708 | 0.1758<br>0.1648 | 0.1724<br>0.1578 | 0.1894 | Ave | | 0.1757 | | | 0.0100 | 7.9 | | 20.0 | | | |
| n-Octadecane | 0.5786 | 0.4969<br>0.4950 | 0.5220<br>0.4662 | 0.5373<br>0.4319 | 0.5750 | Ave | | 0.5128 | | | | 9.9 | | 20.0 | | | |
| Pentachlorophenol | 0.1310 | 0.0630<br>0.1277 | 0.0723<br>0.1247 | 0.0870<br>0.1199 | 0.1058 | Lin1 | -0.089 | 0.1258 | | | 0.0500 | | | | 0.9970 | | 0.9900 |
| Phenanthrene | 1.3770<br>1.1812 | 1.2617<br>1.0880 | 1.2072<br>1.0839 | 1.1682<br>1.0482 | 1.1923 | Ave | | 1.1786 | | | 0.7000 | 8.6 | | 20.0 | | | |
| Anthracene | 1.1316<br>1.2010 | 1.1640<br>1.1155 | 1.1522<br>1.1046 | 1.1498<br>1.0751 | 1.1966 | Ave | | 1.1434 | | | 0.7000 | 3.6 | | 20.0 | | | |
| Carbazole | 1.1181 | 0.9855<br>1.0121 | 1.0010<br>0.9736 | 0.9681<br>0.9190 | 0.9888 | Ave | | 0.9958 | | | 0.0100 | 5.7 | | 20.0 | | | |
| Di-n-butyl phthalate | 1.4278 | 1.0543<br>1.2449 | 1.0537<br>1.1894 | 1.0836<br>1.1345 | 1.2603 | Ave | | 1.1811 | | | 0.0100 | 10.9 | | 20.0 | | | |
| Fluoranthene | 1.1437<br>1.2640 | 1.1063<br>1.1301 | 1.1177<br>1.0882 | 1.0757<br>1.0483 | 1.1538 | Ave | | 1.1253 | | | 0.6000 | 5.5 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19  Calibration End Date: 10/30/2015 16:42  Calibration ID: 31170

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Benzidine | 0.6951 | 0.3632<br>0.6772 | 0.4149<br>0.6916 | 0.4087<br>0.6334 | 0.5454 | Lin1 | -0.483 | 0.6732 | | | | | | | 0.9940 | | 0.9900 |
| Pyrene | 1.2053<br>1.1794 | 1.2231<br>1.1038 | 1.2032<br>1.1259 | 1.1539<br>1.0853 | 1.2062 | Ave | | 1.1651 | | | 0.6000 | 4.3 | | 20.0 | | | |
| Butyl benzyl phthalate | 0.5806 | 0.4429<br>0.5620 | 0.4886<br>0.5776 | 0.4908<br>0.5538 | 0.5224 | Ave | | 0.5273 | | | 0.0100 | 9.4 | | 20.0 | | | |
| Bis(2-ethylhexyl) phthalate | 0.7846 | 0.5860<br>0.7590 | 0.6528<br>0.7817 | 0.6396<br>0.7552 | 0.7101 | Ave | | 0.7086 | | | 0.0100 | 10.5 | | 20.0 | | | |
| 3,3'-Dichlorobenzidine | 0.3701 | 0.2827<br>0.3612 | 0.3127<br>0.3755 | 0.3218<br>0.3615 | 0.3555 | Ave | | 0.3426 | | | 0.0100 | 9.6 | | 20.0 | | | |
| Benzo[a]anthracene | 1.1876<br>1.1606 | 1.1259<br>1.0983 | 1.0942<br>1.1316 | 1.0394<br>1.1004 | 1.0862 | Ave | | 1.1138 | | | 0.8000 | 3.9 | | 20.0 | | | |
| Chrysene | 1.2471<br>1.1079 | 1.2562<br>1.0636 | 1.1353<br>1.0780 | 1.0870<br>1.0408 | 1.1277 | Ave | | 1.1271 | | | 0.7000 | 6.8 | | 20.0 | | | |
| Di-n-octyl phthalate | 1.5859 | 0.8921<br>1.6214 | 1.0660<br>1.6195 | 1.1495<br>1.5780 | 1.2347 | Lin1 | -0.537 | 1.6089 | | | 0.0100 | | | | 0.9960 | | 0.9900 |
| Benzo[b]fluoranthene | 0.6162<br>1.2204 | 0.9128<br>1.1527 | 0.9577<br>1.1512 | 1.0001<br>1.2028 | 1.0888 | Ave | | 1.0336 | | | 0.7000 | 18.4 | | 20.0 | | | |
| Benzo[k]fluoranthene | 1.3131<br>1.3193 | 1.3524<br>1.3266 | 1.2758<br>1.3571 | 1.3772<br>1.2656 | 1.4179 | Ave | | 1.3339 | | | 0.7000 | 3.6 | | 20.0 | | | |
| Benzo[a]pyrene | 0.8263<br>1.2313 | 1.0262<br>1.1934 | 1.0302<br>1.1949 | 1.0741<br>1.1922 | 1.1168 | Ave | | 1.0984 | | | 0.7000 | 11.6 | | 20.0 | | | |
| Indeno[1,2,3-cd]pyrene | 0.7163<br>1.2561 | 0.9705<br>1.2084 | 0.9906<br>1.2361 | 1.1154<br>1.2234 | 1.2408 | Ave | | 1.1064 | | | 0.5000 | 16.5 | | 20.0 | | | |
| Dibenz(a,h)anthracene | 0.5695<br>1.0470 | 0.7363<br>0.9974 | 0.7350<br>1.0228 | 0.8353<br>0.9987 | 0.9400 | Ave | | 0.8758 | | | 0.4000 | 18.9 | | 20.0 | | | |
| Benzo[g,h,i]perylene | 0.8647<br>1.0804 | 0.9477<br>1.0159 | 0.9127<br>1.0424 | 0.9792<br>1.0274 | 1.0726 | Ave | | 0.9937 | | | 0.5000 | 7.4 | | 20.0 | | | |
| 2-Fluorophenol (Surr) | 1.5911 | 1.4167<br>1.5027 | 1.4944<br>1.5649 | 1.5041<br>1.4888 | 1.5108 | Ave | | 1.5092 | | | | 3.5 | | 20.0 | | | |
| Phenol-d5 (Surr) | 2.0299 | 1.9547<br>1.8921 | 2.0082<br>1.9505 | 2.0039<br>1.8641 | 1.9846 | Ave | | 1.9610 | | | | 3.0 | | 20.0 | | | |
| Nitrobenzene-d5 (Surr) | 0.5016<br>0.4763 | 0.4486<br>0.4608 | 0.4786<br>0.4646 | 0.4647<br>0.4426 | 0.4825 | Ave | | 0.4689 | | | | 3.9 | | 20.0 | | | |
| 2-Fluorobiphenyl (Surr) | 1.4656<br>1.2301 | 1.3221<br>1.1942 | 1.2848<br>1.1497 | 1.3000<br>1.1252 | 1.2728 | Ave | | 1.2605 | | | | 8.1 | | 20.0 | | | |
| 2,4,6-Tribromophenol (Surr) | 0.1475 | 0.1223<br>0.1518 | 0.1207<br>0.1515 | 0.1348<br>0.1508 | 0.1415 | Ave | | 0.1401 | | | | 9.2 | | 20.0 | | | |
| Terphenyl-d14 (Surr) | 0.8153<br>0.7449 | 0.7288<br>0.7126 | 0.7158<br>0.7256 | 0.7120<br>0.7014 | 0.7254 | Ave | | 0.7313 | | | | 4.6 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-204569/6 | 51030006.D |
| Level 2 | STD2 240-204569/5 | 51030005.D |
| Level 3 | STD3 240-204569/4 | 51030004.D |
| Level 4 | STD4 240-204569/3 | 51030003.D |
| Level 5 | STD5 240-204569/2 | 51030002.D |
| Level 6 | STD6 240-204569/10 | 51030010.D |
| Level 7 | STD7 240-204569/9 | 51030009.D |
| Level 8 | STD8 240-204569/8 | 51030008.D |
| Level 9 | STD9 240-204569/7 | 51030007.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1,4-Dioxane | DCB | Ave | 494720 | 19536<br>858757 | 45101<br>1002207 | 78767<br>1179056 | 175472 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| N-Nitrosodimethylamine | DCB | Ave | 727704 | 30517<br>1251480 | 56167<br>1478705 | 115169<br>1720943 | 250359 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Pyridine | DCB | Ave | 1312692 | 58896<br>2213458 | 106918<br>2657332 | 210760<br>3089709 | 476330 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzaldehyde | DCB | Ave | 1723444 | 91685<br>2685070 | 177371<br>3127261 | 317074<br>+++++ | 682758 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>+++++ | 10.0 |
| Phenol | DCB | Ave | 1372080 | 63332<br>2238592 | 118575<br>2635317 | 222038<br>3069128 | 498564 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Aniline | DCB | Ave | 1712782 | 76506<br>2822167 | 148036<br>3352650 | 275124<br>3961894 | 604786 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Bis(2-chloroethyl)ether | DCB | Ave | 1165184 | 55798<br>1883482 | 104386<br>2225987 | 193250<br>2628356 | 423165 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Chlorophenol | DCB | Ave | 1003857 | 42048<br>1668277 | 87538<br>1954843 | 158057<br>2306024 | 350205 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| n-Decane | DCB | Ave | 1119398 | 54561<br>1822068 | 107256<br>2096772 | 191841<br>2454439 | 419214 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,3-Dichlorobenzene | DCB | Ave | 1009832 | 45462<br>1668870 | 94475<br>1973407 | 169452<br>2320262 | 365782 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,4-Dichlorobenzene | DCB | Ave | 1025527 | 49389<br>1693331 | 90973<br>2009125 | 172096<br>2369533 | 371977 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzyl alcohol | DCB | Ave | 700878 | 25624<br>1200789 | 52023<br>1423098 | 103687<br>1687761 | 241196 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2-Dichlorobenzene | DCB | Ave | 971991 | 48523<br>1612863 | 88066<br>1929773 | 165378<br>2258088 | 359358 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Methylphenol | DCB | Ave | 1031858 | 45266<br>1698523 | 89970<br>2035078 | 159890<br>2387471 | 365511 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| bis (2-chloroisopropyl) ether | DCB | Ave | 1297319 | 62059<br>2124999 | 121804<br>2502554 | 214776<br>2906661 | 472227 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Indene | DCB | Ave | 3425065 | 165574<br>5482764 | 328167<br>6429527 | 600362<br>7467270 | 1282574 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 3 & 4 Methylphenol | DCB | Ave | 1058372 | 46991<br>1775344 | 86836<br>2093566 | 174167<br>2434286 | 381951 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| N-Nitrosodi-n-propylamine | DCB | Ave | 752762 | 32563<br>1232154 | 65716<br>1447920 | 120161<br>1673167 | 268296 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Acetophenone | DCB | Ave | 1496318 | 71901<br>2512010 | 132596<br>2963572 | 248987<br>3517556 | 555447 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachloroethane | DCB | Ave | 430693 | 20108<br>726077 | 37282<br>860465 | 69054<br>1026455 | 153202 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Nitrobenzene | NPT | Ave | 1232535 | 52171<br>2048374 | 100394<br>2449663 | 197853<br>2964558 | 440896 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Isophorone | NPT | Ave | 2103531 | 86826<br>3593093 | 177542<br>4220065 | 331665<br>4994533 | 753179 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Nitrophenol | NPT | Ave | 491115 | 15975<br>847449 | 33482<br>989517 | 67890<br>1186326 | 168486 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dimethylphenol | NPT | Ave | 1039737 | 46374<br>1762498 | 91248<br>2058849 | 163677<br>2422628 | 380224 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzoic acid | NPT | Lin1 | 1142817 | +++++<br>2058098 | 38740<br>2498369 | 96930<br>2950073 | 285624 | 20.0 | +++++<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Bis(2-chloroethoxy)methane | NPT | Ave | 1244682 | 56184<br>2019878 | 110198<br>2388429 | 201525<br>2812442 | 452022 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dichlorophenol | NPT | Ave | 694975 | 27014<br>1188119 | 58010<br>1406299 | 105633<br>1691338 | 243177 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2,4-Trichlorobenzene | NPT | Ave | 780247 | 35565<br>1315136 | 70146<br>1554140 | 127842<br>1875881 | 288287 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Naphthalene | NPT | Ave | 30876<br>2887186 | 135436<br>4803088 | 258393<br>5581735 | 472164<br>6624747 | 1038816 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Chloroaniline | NPT | Ave | 1318658 | 54609<br>2228978 | 106325<br>2623952 | 193137<br>3120285 | 435662 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,6-Dichlorophenol | NPT | Ave | 703637 | 30354<br>1185386 | 60242<br>1377629 | 106196<br>1675535 | 244055 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorobutadiene | NPT | Ave | 414688 | 20438<br>699720 | 36299<br>823607 | 67299<br>988735 | 147804 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Caprolactam | NPT | Ave | 644498 | 21257<br>1119070 | 48182<br>1324265 | 96949<br>1601386 | 215782 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 4-Chloro-3-methylphenol | NPT | Ave | 914986 | 33914<br>1557755 | 67831<br>1856825 | 130984<br>2233302 | 310185 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Methylnaphthalene | NPT | Ave | 15766<br>1840229 | 87848<br>3059284 | 162653<br>3541203 | 296237<br>4184178 | 663057 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1-Methylnaphthalene | NPT | Ave | 15283<br>1630843 | 72788<br>2686399 | 141746<br>3131422 | 265761<br>3776979 | 581313 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorocyclopentadiene | ANT | Ave | <br>394044 | 11364<br>701729 | 24508<br>841298 | 49561<br>1008324 | 123507 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2,4,5-Tetrachlorobenzene | ANT | Ave | <br>671001 | 30988<br>1118594 | 60915<br>1285990 | 110930<br>1535750 | 239595 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,6-Trichlorophenol | ANT | Ave | <br>489154 | 17900<br>838872 | 37692<br>980224 | 71064<br>1181227 | 160076 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,5-Trichlorophenol | ANT | Ave | <br>501810 | 18213<br>870614 | 41236<br>1011498 | 70318<br>1235185 | 172716 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,1'-Biphenyl | ANT | Ave | 19724<br>2091779 | 95650<br>3462991 | 183903<br>4026917 | 331864<br>4824900 | 728318 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Chloronaphthalene | ANT | Ave | 14887<br>1569866 | 70286<br>2616293 | 137533<br>3089513 | 250588<br>3649714 | 555820 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Nitroaniline | ANT | Ave | <br>614741 | 18048<br>1061375 | 40558<br>1232281 | 83742<br>1478539 | 197887 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dimethyl phthalate | ANT | Ave | <br>1702173 | 72930<br>2843860 | 145058<br>3254556 | 262568<br>3896026 | 588480 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,3-Dinitrobenzene | ANT | Ave | <br>298485 | 7928<br>527703 | 16552<br>607955 | 36557<br>747972 | 93930 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,6-Dinitrotoluene | ANT | Ave | <br>419004 | 14062<br>710781 | 31896<br>815429 | 57470<br>992797 | 135610 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Acenaphthylene | ANT | Ave | 24100<br>2666569 | 105956<br>4418236 | 222529<br>5112162 | 402438<br>6108183 | 909132 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 3-Nitroaniline | ANT | Ave | <br>511901 | 16275<br>874019 | 38619<br>1025857 | 71914<br>1259119 | 165953 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dinitrophenol | ANT | Lin | <br>340757 | +++++<br>711943 | 4284<br>860610 | 12410<br>1055304 | 64957 | <br>20.0 | +++++<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Acenaphthene | ANT | Ave | 16814<br>1656128 | 73639<br>2742805 | 143873<br>3137970 | 262136<br>3799965 | 582412 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Nitrophenol | ANT | Ave | <br>656716 | 15091<br>1126971 | 40091<br>1328703 | 83844<br>1586727 | 196469 | <br>20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 2,4-Dinitrotoluene | ANT | Ave | <br>530683 | 14830<br>929601 | 37139<br>1128606 | 70558<br>1357003 | 177879 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dibenzofuran | ANT | Ave | 22278<br>2207518 | 100992<br>3685353 | 195779<br>4268396 | 347613<br>5091440 | 774124 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,3,4,6-Tetrachlorophenol | ANT | Ave | <br>400411 | 11644<br>657634 | 27986<br>775489 | 56691<br>950049 | 132098 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexadecane | ANT | Ave | <br>1170397 | 53266<br>1772939 | 115676<br>1985743 | 206004<br>+++++ | 437525 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>+++++ | 5.00 |
| Diethyl phthalate | ANT | Ave | <br>1588885 | 73898<br>2662057 | 153677<br>3084203 | 280040<br>3739183 | 604505 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 4-Chlorophenyl phenyl ether | ANT | Ave | 818813 | 35249<br>1392937 | 72831<br>1597010 | 128218<br>1894710 | 280312 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Nitroaniline | ANT | Ave | 536627 | 15768<br>929211 | 36705<br>1058122 | 72711<br>1218830 | 160704 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Fluorene | ANT | Ave | 16049<br>1806392 | 78140<br>2952947 | 153249<br>3365070 | 279644<br>3916619 | 617655 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4,6-Dinitro-2-methylphenol | PHN | Lin1 | 540202 | +++++<br>1039361 | 18320<br>1201923 | 48290<br>1468698 | 142860 | 20.0 | +++++<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| N-Nitrosodiphenylamine | PHN | Ave | 2412476 | 107217<br>3871646 | 219238<br>4466925 | 395911<br>5367714 | 864865 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | PHN | Ave | 2408330 | 96968<br>3930371 | 197817<br>4489206 | 372021<br>5272898 | 839206 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Azobenzene | PHN | Ave | 2408283 | 97367<br>3930767 | 197585<br>4489447 | 371861<br>5274047 | 839582 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Bromophenyl phenyl ether | PHN | Ave | 440500 | 17809<br>709837 | 38709<br>840209 | 67242<br>995151 | 144527 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorobenzene | PHN | Ave | 4393<br>427015 | 19589<br>690109 | 38717<br>804235 | 68470<br>958448 | 141862 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Atrazine | PHN | Ave | 855420 | 33504<br>1352914 | 70696<br>1545180 | 125346<br>1842351 | 286069 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| n-Octadecane | PHN | Ave | 1227360 | 48071<br>1960495 | 104943<br>2185258 | 195341<br>2521202 | 434198 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Pentachlorophenol | PHN | Lin1 | 555699 | 12198<br>1011380 | 29080<br>1169285 | 63227<br>1399974 | 159821 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Phenanthrene | PHN | Ave | 26012<br>2505761 | 122072<br>4309272 | 242690<br>5081149 | 424735<br>6119446 | 900393 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Anthracene | PHN | Ave | 21376<br>2547843 | 112620<br>4418412 | 231620<br>5178390 | 418036<br>6276450 | 903640 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Carbazole | PHN | Ave | 2371890 | 95351<br>4008777 | 201234<br>4564046 | 351979<br>5364842 | 746715 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Di-n-butyl phthalate | PHN | Ave | 3028900 | 102002<br>4930803 | 211817<br>5575661 | 393979<br>6623099 | 951717 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Fluoranthene | PHN | Ave | 21605<br>2681466 | 107034<br>4476252 | 224697<br>5101161 | 391101<br>6120088 | 871324 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzidine | CRY | Lin1 | 3263077 | 68262<br>5801270 | 165601<br>6638385 | 293457<br>7576128 | 855611 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Pyrene | CRY | Ave | 21112<br>2768391 | 114934<br>4728230 | 240149<br>5403729 | 414229<br>6491243 | 946189 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Butyl benzyl phthalate | CRY | Ave | 1362886 | 41616<br>2407141 | 97515<br>2772322 | 176174<br>3312554 | 409808 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Bis(2-ethylhexyl) phthalate | CRY | Ave | 1841586 | 55068<br>3251044 | 130297<br>3751555 | 229614<br>4516996 | 557021 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 3,3'-Dichlorobenzidine | CRY | Ave | 1737283 | 53130<br>3094408 | 124821<br>3604860 | 231023<br>4323701 | 557748 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Benzo[a]anthracene | CRY | Ave | 20802<br>2724169 | 105800<br>4704482 | 218392<br>5431316 | 373133<br>6581213 | 852065 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Chrysene | CRY | Ave | 21844<br>2600462 | 118051<br>4555837 | 226582<br>5173767 | 390188<br>6225007 | 884584 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Di-n-octyl phthalate | PRY | Lin1 | 3164188 | 72150<br>5868029 | 178621<br>6734392 | 343630<br>8162171 | 780776 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[b]fluoranthene | PRY | Ave | 9374<br>2434889 | 73820<br>4171749 | 160463<br>4786938 | 298966<br>6221630 | 688506 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[k]fluoranthene | PRY | Ave | 19974<br>2632345 | 109374<br>4801140 | 213764<br>5643347 | 411682<br>6546419 | 896627 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[a]pyrene | PRY | Ave | 12569<br>2456604 | 82995<br>4319074 | 172613<br>4968958 | 321073<br>6166812 | 706216 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Indeno[1,2,3-cd]pyrene | PRY | Ave | 10896<br>2506233 | 78485<br>4373399 | 165972<br>5139962 | 333419<br>6327726 | 784631 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dibenz(a,h)anthracene | PRY | Ave | 8663<br>2089005 | 59548<br>3609922 | 123160<br>4253007 | 249705<br>5165905 | 594408 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[g,h,i]perylene | PRY | Ave | 13154<br>2155633 | 76647<br>3676705 | 152929<br>4334830 | 292707<br>5314104 | 678278 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Fluorophenol (Surr) | DCB | Ave | 998997 | 41294<br>1670627 | 84646<br>1982372 | 157568<br>2316746 | 349930 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Phenol-d5 (Surr) | DCB | Ave | 1274518 | 56976<br>2103557 | 113748<br>2470874 | 209937<br>2900830 | 459672 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Nitrobenzene-d5 (Surr) | NPT | Ave | 11780<br>1279836 | 52185<br>2181131 | 108855<br>2622430 | 199480<br>3089135 | 456387 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Fluorobiphenyl (Surr) | ANT | Ave | 16631<br>1709206 | 73911<br>2853677 | 149335<br>3309526 | 273305<br>3962722 | 599773 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,6-Tribromophenol (Surr) | ANT | Ave | 204995 | 6837<br>362734 | 14034<br>436170 | 28334<br>531122 | 66695 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Terphenyl-d14 (Surr) | CRY | Ave | 14281<br>1748429 | 68485<br>3052200 | 142868<br>3482351 | 255585<br>4195262 | 569027 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin = Linear ISTD |
| Lin1 = Linear 1/conc ISTD |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID (1): CCV 240-206233/2 CCVIS Instrument ID (1): A4HP9

GC Column (1): RXI-5SILMS ID: 0.45(mm) Date Analyzed (1): 11/11/2015 07:49

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 2-Fluorophenol (Surr) | 5.04 | 100.0 |
| Benzaldehyde | 5.89 | 100.0 |
| Phenol-d5 (Surr) | 5.95 | 100.0 |
| Phenol | 5.96 | 100.0 |
| Bis(2-chloroethyl)ether | 6.03 | 100.0 |
| 2-Chlorophenol | 6.12 | 100.0 |
| 2-Methylphenol | 6.50 | 100.0 |
| bis (2-chloroisopropyl) ether | 6.52 | 100.0 |
| N-Nitrosodi-n-propylamine | 6.63 | 100.0 |
| 3 & 4 Methylphenol | 6.63 | 100.0 |
| Acetophenone | 6.64 | 100.0 |
| Hexachloroethane | 6.76 | 100.0 |
| Nitrobenzene-d5 (Surr) | 6.78 | 100.0 |
| Nitrobenzene | 6.80 | 100.0 |
| Isophorone | 7.00 | 100.0 |
| 2-Nitrophenol | 7.08 | 100.0 |
| 2,4-Dimethylphenol | 7.09 | 100.0 |
| Bis(2-chloroethoxy)methane | 7.16 | 100.0 |
| 2,4-Dichlorophenol | 7.29 | 100.0 |
| Naphthalene | 7.44 | 100.0 |
| 4-Chloroaniline | 7.46 | 100.0 |
| Hexachlorobutadiene | 7.54 | 100.0 |
| Caprolactam | 7.75 | 100.0 |
| 4-Chloro-3-methylphenol | 7.86 | 100.0 |
| 2-Methylnaphthalene | 8.02 | 100.0 |
| Hexachlorocyclopentadiene | 8.16 | 100.0 |
| 2,4,6-Trichlorophenol | 8.26 | 100.0 |
| 2,4,5-Trichlorophenol | 8.30 | 100.0 |
| 2-Fluorobiphenyl (Surr) | 8.32 | 100.0 |
| 1,1'-Biphenyl | 8.41 | 100.0 |
| 2-Chloronaphthalene | 8.45 | 100.0 |
| 2-Nitroaniline | 8.51 | 100.0 |
| Dimethyl phthalate | 8.64 | 100.0 |
| 2,6-Dinitrotoluene | 8.70 | 100.0 |
| Acenaphthylene | 8.80 | 100.0 |
| 3-Nitroaniline | 8.86 | 100.0 |
| 2,4-Dinitrophenol | 8.94 | 100.0 |
| Acenaphthene | 8.95 | 100.0 |
| 4-Nitrophenol | 8.98 | 100.0 |
| 2,4-Dinitrotoluene | 9.05 | 100.0 |
| Dibenzofuran | 9.09 | 100.0 |
| Diethyl phthalate | 9.23 | 100.0 |
| 4-Chlorophenyl phenyl ether | 9.36 | 100.0 |
| 4-Nitroaniline | 9.38 | 100.0 |
| Fluorene | 9.39 | 100.0 |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID (1): CCV 240-206233/2 CCVIS Instrument ID (1): A4HP9

GC Column (1): RXI-5SILMS ID: 0.45(mm) Date Analyzed (1): 11/11/2015 07:49

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 4,6-Dinitro-2-methylphenol | 9.40 | 100.0 |
| N-Nitrosodiphenylamine | 9.45 | 100.0 |
| 2,4,6-Tribromophenol (Surr) | 9.59 | 100.0 |
| 4-Bromophenyl phenyl ether | 9.78 | 100.0 |
| Hexachlorobenzene | 9.87 | 100.0 |
| Atrazine | 9.88 | 100.0 |
| Pentachlorophenol | 10.02 | 100.0 |
| Phenanthrene | 10.21 | 100.0 |
| Anthracene | 10.25 | 100.0 |
| Carbazole | 10.37 | 100.0 |
| Di-n-butyl phthalate | 10.60 | 100.0 |
| Fluoranthene | 11.22 | 100.0 |
| Pyrene | 11.42 | 100.0 |
| Terphenyl-d14 (Surr) | 11.51 | 100.0 |
| Butyl benzyl phthalate | 11.88 | 100.0 |
| Bis(2-ethylhexyl) phthalate | 12.35 | 100.0 |
| 3,3'-Dichlorobenzidine | 12.39 | 100.0 |
| Benzo[a]anthracene | 12.45 | 100.0 |
| Chrysene | 12.49 | 100.0 |
| Di-n-octyl phthalate | 12.98 | 100.0 |
| Benzo[b]fluoranthene | 13.65 | 35.70 |
| Benzo[k]fluoranthene | 13.68 | 100.0 |
| Benzo[a]pyrene | 14.10 | 100.0 |
| Indeno[1,2,3-cd]pyrene | 16.06 | 100.0 |
| Dibenz(a,h)anthracene | 16.07 | 100.0 |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID (1): CCV 240-206256/2 CCVIS Instrument ID (1): A4HP10

GC Column (1): DB-624 ID: 0.18(mm) Date Analyzed (1): 11/11/2015 09:34

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 2-Fluorophenol (Surr) | 4.95 | 100.0 |
| Benzaldehyde | 5.76 | 100.0 |
| Phenol-d5 (Surr) | 5.82 | 100.0 |
| Phenol | 5.84 | 100.0 |
| Bis(2-chloroethyl)ether | 5.89 | 100.0 |
| 2-Chlorophenol | 5.98 | 100.0 |
| 2-Methylphenol | 6.36 | 100.0 |
| bis (2-chloroisopropyl) ether | 6.38 | 100.0 |
| N-Nitrosodi-n-propylamine | 6.49 | 100.0 |
| Acetophenone | 6.50 | 100.0 |
| 3 & 4 Methylphenol | 6.50 | 100.0 |
| Hexachloroethane | 6.61 | 100.0 |
| Nitrobenzene-d5 (Surr) | 6.63 | 100.0 |
| Nitrobenzene | 6.65 | 100.0 |
| Isophorone | 6.85 | 100.0 |
| 2-Nitrophenol | 6.93 | 100.0 |
| 2,4-Dimethylphenol | 6.96 | 100.0 |
| Bis(2-chloroethoxy)methane | 7.02 | 100.0 |
| 2,4-Dichlorophenol | 7.15 | 100.0 |
| Naphthalene | 7.29 | 100.0 |
| 4-Chloroaniline | 7.32 | 100.0 |
| Hexachlorobutadiene | 7.39 | 100.0 |
| Caprolactam | 7.61 | 100.0 |
| 4-Chloro-3-methylphenol | 7.73 | 100.0 |
| 2-Methylnaphthalene | 7.88 | 100.0 |
| Hexachlorocyclopentadiene | 8.02 | 100.0 |
| 2,4,6-Trichlorophenol | 8.12 | 100.0 |
| 2,4,5-Trichlorophenol | 8.16 | 100.0 |
| 2-Fluorobiphenyl (Surr) | 8.18 | 100.0 |
| 1,1'-Biphenyl | 8.27 | 100.0 |
| 2-Chloronaphthalene | 8.30 | 100.0 |
| 2-Nitroaniline | 8.37 | 100.0 |
| Dimethyl phthalate | 8.50 | 100.0 |
| 2,6-Dinitrotoluene | 8.56 | 100.0 |
| Acenaphthylene | 8.66 | 100.0 |
| 3-Nitroaniline | 8.71 | 100.0 |
| Acenaphthene | 8.80 | 100.0 |
| 2,4-Dinitrophenol | 8.81 | 100.0 |
| 4-Nitrophenol | 8.86 | 100.0 |
| 2,4-Dinitrotoluene | 8.91 | 100.0 |
| Dibenzofuran | 8.95 | 100.0 |
| Diethyl phthalate | 9.09 | 100.0 |
| 4-Chlorophenyl phenyl ether | 9.21 | 100.0 |
| 4-Nitroaniline | 9.24 | 100.0 |
| Fluorene | 9.24 | 100.0 |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID (1): CCV 240-206256/2 CCVIS Instrument ID (1): A4HP10

GC Column (1): DB-624 ID: 0.18(mm) Date Analyzed (1): 11/11/2015 09:34

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 4,6-Dinitro-2-methylphenol | 9.27 | 100.0 |
| N-Nitrosodiphenylamine | 9.31 | 100.0 |
| 2,4,6-Tribromophenol (Surr) | 9.45 | 100.0 |
| 4-Bromophenyl phenyl ether | 9.63 | 100.0 |
| Hexachlorobenzene | 9.72 | 100.0 |
| Atrazine | 9.74 | 100.0 |
| Pentachlorophenol | 9.88 | 100.0 |
| Phenanthrene | 10.06 | 100.0 |
| Anthracene | 10.11 | 100.0 |
| Carbazole | 10.23 | 100.0 |
| Di-n-butyl phthalate | 10.46 | 100.0 |
| Fluoranthene | 11.09 | 100.0 |
| Pyrene | 11.32 | 100.0 |
| Terphenyl-d14 (Surr) | 11.42 | 100.0 |
| Butyl benzyl phthalate | 11.91 | 100.0 |
| Bis(2-ethylhexyl) phthalate | 12.61 | 100.0 |
| 3,3'-Dichlorobenzidine | 12.62 | 100.0 |
| Benzo[a]anthracene | 12.69 | 100.0 |
| Chrysene | 12.74 | 100.0 |
| Di-n-octyl phthalate | 13.54 | 100.0 |
| Benzo[b]fluoranthene | 14.28 | 25.90 |
| Benzo[k]fluoranthene | 14.33 | 100.0 |
| Benzo[a]pyrene | 14.80 | 100.0 |
| Indeno[1,2,3-cd]pyrene | 16.67 | 100.0 |
| Dibenz(a,h)anthracene | 16.68 | 100.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-200091/11 Calibration Date: 10/02/2015 14:28

Instrument ID: A4HP10 Calib Start Date: 10/02/2015 10:48

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/02/2015 14:04

Lab File ID: 51002011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.6903 | 0.6472 | | 9.38 | 10.0 | -6.2 | 30.0 |
| N-Nitrosodimethylamine | Ave | 0.9253 | 0.7776 | | 8.40 | 10.0 | -16.0 | 30.0 |
| Pyridine | Ave | 1.684 | 1.634 | | 9.70 | 10.0 | -3.0 | 30.0 |
| Benzaldehyde | Ave | 1.111 | 1.071 | 0.0100 | 19.3 | 20.0 | -3.6 | 30.0 |
| Phenol | Ave | 1.891 | 1.832 | 0.8000 | 9.69 | 10.0 | -3.1 | 30.0 |
| Aniline | Ave | 2.402 | 2.297 | | 9.57 | 10.0 | -4.3 | 30.0 |
| Bis(2-chloroethyl)ether | Ave | 1.506 | 1.415 | 0.7000 | 9.40 | 10.0 | -6.0 | 30.0 |
| 2-Chlorophenol | Ave | 1.543 | 1.512 | 0.8000 | 9.80 | 10.0 | -2.0 | 30.0 |
| n-Decane | Ave | 1.911 | 1.813 | | 9.48 | 10.0 | -5.2 | 30.0 |
| 1,3-Dichlorobenzene | Ave | 1.579 | 1.531 | | 9.69 | 10.0 | -3.1 | 30.0 |
| 1,4-Dichlorobenzene | Ave | 1.615 | 1.563 | | 9.68 | 10.0 | -3.2 | 30.0 |
| Benzyl alcohol | Ave | 1.025 | 0.9929 | | 9.69 | 10.0 | -3.1 | 30.0 |
| 1,2-Dichlorobenzene | Ave | 1.531 | 1.463 | | 9.56 | 10.0 | -4.4 | 30.0 |
| 2-Methylphenol | Ave | 1.412 | 1.403 | 0.7000 | 9.93 | 10.0 | -0.7 | 30.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.121 | 2.000 | | 9.43 | 10.0 | -5.7 | 30.0 |
| Indene | Ave | 2.534 | 2.356 | | 18.6 | 20.0 | -7.0 | 30.0 |
| 3 & 4 Methylphenol | Ave | 1.469 | 1.437 | | 9.78 | 10.0 | -2.2 | 30.0 |
| N-Nitrosodi-n-propylamine | Ave | 0.8785 | 0.8794 | 0.5000 | 10.0 | 10.0 | 0.1 | 30.0 |
| Acetophenone | Ave | 1.988 | 1.939 | 0.0100 | 9.76 | 10.0 | -2.4 | 30.0 |
| Hexachloroethane | Ave | 0.6359 | 0.6318 | 0.3000 | 9.94 | 10.0 | -0.6 | 30.0 |
| Nitrobenzene | Ave | 0.3694 | 0.3535 | 0.2000 | 9.57 | 10.0 | -4.3 | 30.0 |
| Isophorone | Ave | 0.6852 | 0.6871 | 0.4000 | 10.0 | 10.0 | 0.3 | 30.0 |
| 2-Nitrophenol | Ave | 0.1669 | 0.1805 | 0.1000 | 10.8 | 10.0 | 8.1 | 30.0 |
| 2,4-Dimethylphenol | Ave | 0.3580 | 0.3438 | 0.2000 | 9.61 | 10.0 | -3.9 | 30.0 |
| Benzoic acid | Lin1 | | 0.2265 | | 21.2 | 20.0 | 5.9 | 30.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4236 | 0.4028 | 0.3000 | 9.51 | 10.0 | -4.9 | 30.0 |
| 2,4-Dichlorophenol | Ave | 0.2940 | 0.2826 | 0.2000 | 9.61 | 10.0 | -3.9 | 30.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.3022 | 0.2949 | | 9.76 | 10.0 | -2.4 | 30.0 |
| Naphthalene | Ave | 1.084 | 1.032 | 0.7000 | 9.52 | 10.0 | -4.8 | 30.0 |
| 4-Chloroaniline | Ave | 0.4756 | 0.4464 | 0.0100 | 9.39 | 10.0 | -6.1 | 30.0 |
| 2,6-Dichlorophenol | Ave | 0.2949 | 0.2868 | | 9.73 | 10.0 | -2.7 | 30.0 |
| Hexachlorobutadiene | Ave | 0.1868 | 0.1827 | 0.0100 | 9.78 | 10.0 | -2.2 | 30.0 |
| Caprolactam | Ave | 0.1164 | 0.1183 | 0.0100 | 20.3 | 20.0 | 1.6 | 30.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3033 | 0.2997 | 0.2000 | 9.88 | 10.0 | -1.2 | 30.0 |
| 2-Methylnaphthalene | Ave | 0.7006 | 0.6481 | 0.4000 | 9.25 | 10.0 | -7.5 | 30.0 |
| 1-Methylnaphthalene | Ave | 0.6254 | 0.6114 | | 9.78 | 10.0 | -2.2 | 30.0 |
| Hexachlorocyclopentadiene | Ave | 0.2689 | 0.2968 | 0.0500 | 11.0 | 10.0 | 10.4 | 30.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4998 | 0.4717 | 0.0100 | 9.44 | 10.0 | -5.6 | 30.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3619 | 0.3701 | 0.2000 | 10.2 | 10.0 | 2.3 | 30.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3794 | 0.3824 | 0.2000 | 10.1 | 10.0 | 0.8 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-200091/11 Calibration Date: 10/02/2015 14:28

Instrument ID: A4HP10 Calib Start Date: 10/02/2015 10:48

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/02/2015 14:04

Lab File ID: 51002011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.409 | 1.323 | 0.0100 | 9.39 | 10.0 | -6.1 | 30.0 |
| 2-Chloronaphthalene | Ave | 1.182 | 1.109 | 0.8000 | 9.38 | 10.0 | -6.2 | 30.0 |
| 2-Nitroaniline | Ave | 0.3503 | 0.3633 | 0.0100 | 10.4 | 10.0 | 3.7 | 30.0 |
| Dimethyl phthalate | Ave | 1.285 | 1.249 | 0.0100 | 9.72 | 10.0 | -2.8 | 30.0 |
| 1,3-Dinitrobenzene | Lin1 | | 0.2089 | | 10.0 | 10.0 | 0.3 | 30.0 |
| 2,6-Dinitrotoluene | Ave | 0.2818 | 0.2952 | | 10.5 | 10.0 | 4.8 | 30.0 |
| Acenaphthylene | Ave | 1.763 | 1.642 | 0.9000 | 9.31 | 10.0 | -6.9 | 30.0 |
| 3-Nitroaniline | Ave | 0.3239 | 0.3357 | 0.0100 | 10.4 | 10.0 | 3.6 | 30.0 |
| 2,4-Dinitrophenol | Lin1 | | 0.1483 | 0.0100 | 20.9 | 20.0 | 4.7 | 30.0 |
| Acenaphthene | Ave | 1.075 | 1.012 | 0.9000 | 9.41 | 10.0 | -5.9 | 30.0 |
| 4-Nitrophenol | Ave | 0.1702 | 0.1772 | | 20.8 | 20.0 | 4.1 | 30.0 |
| 2,4-Dinitrotoluene | Ave | 0.3666 | 0.3977 | 0.2000 | 10.8 | 10.0 | 8.5 | 30.0 |
| Dibenzofuran | Ave | 1.623 | 1.525 | 0.8000 | 9.39 | 10.0 | -6.1 | 30.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.3179 | 0.3218 | 0.0100 | 10.1 | 10.0 | 1.2 | 30.0 |
| Diethyl phthalate | Ave | 1.172 | 1.149 | 0.0100 | 9.80 | 10.0 | -2.0 | 30.0 |
| Hexadecane | Ave | 0.5546 | 0.5084 | | 9.17 | 10.0 | -8.3 | 30.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.6110 | 0.5884 | 0.4000 | 9.63 | 10.0 | -3.7 | 30.0 |
| 4-Nitroaniline | Ave | 0.3149 | 0.3341 | 0.0100 | 10.6 | 10.0 | 6.1 | 30.0 |
| Fluorene | Ave | 1.322 | 1.271 | 0.9000 | 9.61 | 10.0 | -3.9 | 30.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1420 | 0.0100 | 20.8 | 20.0 | 3.8 | 30.0 |
| N-Nitrosodiphenylamine | Ave | 0.5710 | 0.5568 | 0.0100 | 19.5 | 20.0 | -2.5 | 30.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 0.7994 | 0.8227 | | 10.3 | 10.0 | 2.9 | 30.0 |
| Azobenzene | Ave | 0.7997 | 0.8228 | | 10.3 | 10.0 | 2.9 | 30.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.2298 | 0.2289 | 0.1000 | 9.96 | 10.0 | -0.4 | 30.0 |
| Hexachlorobenzene | Ave | 0.2582 | 0.2429 | 0.1000 | 9.41 | 10.0 | -5.9 | 30.0 |
| Atrazine | Ave | 0.2140 | 0.2028 | 0.0100 | 19.0 | 20.0 | -5.2 | 30.0 |
| n-Octadecane | Qua | | 0.7442 | | 9.66 | 10.0 | -3.4 | 30.0 |
| Pentachlorophenol | Ave | 0.1432 | 0.1612 | 0.0500 | 22.5 | 20.0 | 12.5 | 30.0 |
| Phenanthrene | Ave | 1.177 | 1.107 | 0.7000 | 9.41 | 10.0 | -5.9 | 30.0 |
| Anthracene | Ave | 1.190 | 1.153 | 0.7000 | 9.69 | 10.0 | -3.1 | 30.0 |
| Carbazole | Ave | 1.127 | 1.062 | 0.0100 | 9.42 | 10.0 | -5.8 | 30.0 |
| Di-n-butyl phthalate | Ave | 1.380 | 1.370 | 0.0100 | 9.93 | 10.0 | -0.7 | 30.0 |
| Fluoranthene | Ave | 1.321 | 1.260 | 0.6000 | 9.54 | 10.0 | -4.6 | 30.0 |
| Benzidine | Ave | 0.5843 | 0.5955 | | 20.4 | 20.0 | 1.9 | 30.0 |
| Pyrene | Ave | 1.135 | 1.058 | 0.6000 | 9.33 | 10.0 | -6.7 | 30.0 |
| Butyl benzyl phthalate | Ave | 0.5142 | 0.5113 | 0.0100 | 9.94 | 10.0 | -0.6 | 30.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3952 | 0.3739 | 0.0100 | 18.9 | 20.0 | -5.4 | 30.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.6579 | 0.6437 | 0.0100 | 9.78 | 10.0 | -2.2 | 30.0 |
| Benzo[a]anthracene | Ave | 1.126 | 1.090 | 0.8000 | 9.68 | 10.0 | -3.2 | 30.0 |
| Chrysene | Ave | 1.076 | 1.070 | 0.7000 | 9.95 | 10.0 | -0.5 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-200091/11 Calibration Date: 10/02/2015 14:28

Instrument ID: A4HP10 Calib Start Date: 10/02/2015 10:48

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/02/2015 14:04

Lab File ID: 51002011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Ave | 1.371 | 1.402 | 0.0100 | 10.2 | 10.0 | 2.3 | 30.0 |
| Benzo[b]fluoranthene | Ave | 1.128 | 1.130 | 0.7000 | 10.0 | 10.0 | 0.2 | 30.0 |
| Benzo[k]fluoranthene | Ave | 1.274 | 1.168 | 0.7000 | 9.17 | 10.0 | -8.3 | 30.0 |
| Benzo[a]pyrene | Ave | 1.120 | 1.096 | 0.7000 | 9.78 | 10.0 | -2.2 | 30.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.233 | 1.234 | 0.5000 | 10.0 | 10.0 | 0.0 | 30.0 |
| Dibenz(a,h)anthracene | Ave | 1.024 | 1.069 | 0.4000 | 10.4 | 10.0 | 4.4 | 30.0 |
| Benzo[g,h,i]perylene | Ave | 1.092 | 1.107 | 0.5000 | 10.1 | 10.0 | 1.3 | 30.0 |
| 2-Fluorophenol (Surr) | Ave | 1.486 | 1.530 | | 10.3 | 10.0 | 3.0 | 30.0 |
| Phenol-d5 (Surr) | Ave | 1.814 | 1.897 | | 10.5 | 10.0 | 4.6 | 30.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.3982 | 0.4088 | | 10.3 | 10.0 | 2.6 | 30.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.297 | 1.284 | | 9.90 | 10.0 | -1.0 | 30.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1891 | 0.2114 | | 11.2 | 10.0 | 11.8 | 30.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7619 | 0.7747 | | 10.2 | 10.0 | 1.7 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206256/2 Calibration Date: 11/11/2015 09:34

Instrument ID: A4HP10 Calib Start Date: 10/02/2015 10:48

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/02/2015 14:04

Lab File ID: 51111002.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.6903 | 0.5705 | | 8.27 | 10.0 | -17.3 | 20.0 |
| N-Nitrosodimethylamine | Ave | 0.9253 | 0.7500 | | 8.11 | 10.0 | -18.9 | 20.0 |
| Pyridine | Ave | 1.684 | 1.387 | | 8.23 | 10.0 | -17.7 | 20.0 |
| Benzaldehyde | Ave | 1.111 | 0.6352 | 0.0100 | 11.4 | 20.0 | -42.8* | 20.0 |
| Phenol | Ave | 1.891 | 1.893 | 0.8000 | 10.0 | 10.0 | 0.0 | 20.0 |
| Aniline | Ave | 2.402 | 2.168 | | 9.03 | 10.0 | -9.7 | 20.0 |
| Bis(2-chloroethyl)ether | Ave | 1.506 | 1.420 | 0.7000 | 9.43 | 10.0 | -5.7 | 20.0 |
| 2-Chlorophenol | Ave | 1.543 | 1.438 | 0.8000 | 9.32 | 10.0 | -6.8 | 20.0 |
| n-Decane | Ave | 1.911 | 1.929 | | 10.1 | 10.0 | 0.9 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.579 | 1.537 | | 9.73 | 10.0 | -2.7 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.615 | 1.575 | | 9.76 | 10.0 | -2.4 | 20.0 |
| Benzyl alcohol | Ave | 1.025 | 0.3563 | | 3.48 | 10.0 | -65.2* | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.531 | 1.486 | | 9.70 | 10.0 | -3.0 | 20.0 |
| 2-Methylphenol | Ave | 1.412 | 1.308 | 0.7000 | 9.27 | 10.0 | -7.3 | 20.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.121 | 2.250 | | 10.6 | 10.0 | 6.1 | 20.0 |
| Indene | Ave | 2.534 | 2.469 | | 19.5 | 20.0 | -2.6 | 20.0 |
| N-Nitrosodi-n-propylamine | Ave | 0.8785 | 0.8567 | 0.5000 | 9.75 | 10.0 | -2.5 | 20.0 |
| 3 & 4 Methylphenol | Ave | 1.469 | 1.380 | | 9.39 | 10.0 | -6.1 | 20.0 |
| Acetophenone | Ave | 1.988 | 1.975 | 0.0100 | 9.93 | 10.0 | -0.7 | 20.0 |
| Hexachloroethane | Ave | 0.6359 | 0.6671 | 0.3000 | 10.5 | 10.0 | 4.9 | 20.0 |
| Nitrobenzene | Ave | 0.3694 | 0.3589 | 0.2000 | 9.72 | 10.0 | -2.8 | 20.0 |
| Isophorone | Ave | 0.6852 | 0.6730 | 0.4000 | 9.82 | 10.0 | -1.8 | 20.0 |
| 2-Nitrophenol | Ave | 0.1669 | 0.1844 | 0.1000 | 11.0 | 10.0 | 10.5 | 20.0 |
| 2,4-Dimethylphenol | Ave | 0.3580 | 0.3249 | 0.2000 | 9.07 | 10.0 | -9.3 | 20.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4236 | 0.4023 | 0.3000 | 9.50 | 10.0 | -5.0 | 20.0 |
| Benzoic acid | Lin1 | | 0.1787 | | 17.0 | 20.0 | -15.1 | 20.0 |
| 2,4-Dichlorophenol | Ave | 0.2940 | 0.2941 | 0.2000 | 10.0 | 10.0 | 0.0 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.3022 | 0.3090 | | 10.2 | 10.0 | 2.2 | 20.0 |
| Naphthalene | Ave | 1.084 | 1.060 | 0.7000 | 9.78 | 10.0 | -2.2 | 20.0 |
| 4-Chloroaniline | Ave | 0.4756 | 0.4439 | 0.0100 | 9.33 | 10.0 | -6.7 | 20.0 |
| 2,6-Dichlorophenol | Ave | 0.2949 | 0.2816 | | 9.55 | 10.0 | -4.5 | 20.0 |
| Hexachlorobutadiene | Ave | 0.1868 | 0.1936 | 0.0100 | 10.4 | 10.0 | 3.6 | 20.0 |
| Caprolactam | Ave | 0.1164 | 0.1013 | 0.0100 | 17.4 | 20.0 | -12.9 | 20.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3033 | 0.3014 | 0.2000 | 9.94 | 10.0 | -0.6 | 20.0 |
| 2-Methylnaphthalene | Ave | 0.7006 | 0.6945 | 0.4000 | 9.91 | 10.0 | -0.9 | 20.0 |
| 1-Methylnaphthalene | Ave | 0.6254 | 0.6095 | | 9.75 | 10.0 | -2.5 | 20.0 |
| Hexachlorocyclopentadiene | Ave | 0.2689 | 0.1923 | 0.0500 | 7.15 | 10.0 | -28.5* | 20.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4998 | 0.5133 | 0.0100 | 10.3 | 10.0 | 2.7 | 20.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3619 | 0.3550 | 0.2000 | 9.81 | 10.0 | -1.9 | 20.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3794 | 0.3767 | 0.2000 | 9.93 | 10.0 | -0.7 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206256/2 | Calibration Date: 11/11/2015 09:34

Instrument ID: A4HP10 | Calib Start Date: 10/02/2015 10:48

GC Column: DB-624 ID: 0.18(mm) | Calib End Date: 10/02/2015 14:04

Lab File ID: 51111002.D | Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.409 | 1.370 | 0.0100 | 9.72 | 10.0 | -2.8 | 20.0 |
| 2-Chloronaphthalene | Ave | 1.182 | 1.123 | 0.8000 | 9.50 | 10.0 | -5.0 | 20.0 |
| 2-Nitroaniline | Ave | 0.3503 | 0.3581 | 0.0100 | 10.2 | 10.0 | 2.2 | 20.0 |
| Dimethyl phthalate | Ave | 1.285 | 1.285 | 0.0100 | 10.0 | 10.0 | 0.0 | 20.0 |
| 1,3-Dinitrobenzene | Lin1 | | 0.2007 | | 9.65 | 10.0 | -3.5 | 20.0 |
| 2,6-Dinitrotoluene | Ave | 0.2818 | 0.2901 | | 10.3 | 10.0 | 3.0 | 20.0 |
| Acenaphthylene | Ave | 1.763 | 1.739 | 0.9000 | 9.87 | 10.0 | -1.3 | 20.0 |
| 3-Nitroaniline | Ave | 0.3239 | 0.3137 | 0.0100 | 9.68 | 10.0 | -3.2 | 20.0 |
| Acenaphthene | Ave | 1.075 | 1.050 | 0.9000 | 9.77 | 10.0 | -2.3 | 20.0 |
| 2,4-Dinitrophenol | Lin1 | | 0.1133 | 0.0100 | 16.5 | 20.0 | -17.5 | 20.0 |
| 4-Nitrophenol | Ave | 0.1702 | 0.1831 | | 21.5 | 20.0 | 7.6 | 20.0 |
| 2,4-Dinitrotoluene | Ave | 0.3666 | 0.3849 | 0.2000 | 10.5 | 10.0 | 5.0 | 20.0 |
| Dibenzofuran | Ave | 1.623 | 1.575 | 0.8000 | 9.70 | 10.0 | -3.0 | 20.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.3179 | 0.2913 | 0.0100 | 9.16 | 10.0 | -8.4 | 20.0 |
| Diethyl phthalate | Ave | 1.172 | 1.234 | 0.0100 | 10.5 | 10.0 | 5.3 | 20.0 |
| Hexadecane | Ave | 0.5546 | 0.5740 | | 10.3 | 10.0 | 3.5 | 20.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.6110 | 0.5883 | 0.4000 | 9.63 | 10.0 | -3.7 | 20.0 |
| 4-Nitroaniline | Ave | 0.3149 | 0.2932 | 0.0100 | 9.31 | 10.0 | -6.9 | 20.0 |
| Fluorene | Ave | 1.322 | 1.299 | 0.9000 | 9.82 | 10.0 | -1.8 | 20.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1067 | 0.0100 | 16.0 | 20.0 | -20.0 | 20.0 |
| N-Nitrosodiphenylamine | Ave | 0.5710 | 0.5894 | 0.0100 | 20.6 | 20.0 | 3.2 | 20.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 0.7994 | 0.7848 | | 9.81 | 10.0 | -1.8 | 20.0 |
| Azobenzene | Ave | 0.7997 | 0.7841 | | 9.80 | 10.0 | -2.0 | 20.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.2298 | 0.2307 | 0.1000 | 10.0 | 10.0 | 0.4 | 20.0 |
| Hexachlorobenzene | Ave | 0.2582 | 0.2467 | 0.1000 | 9.55 | 10.0 | -4.5 | 20.0 |
| Atrazine | Ave | 0.2140 | 0.2111 | 0.0100 | 19.7 | 20.0 | -1.4 | 20.0 |
| n-Octadecane | Qua | | 0.9708 | | 13.8 | 10.0 | 37.9* | 20.0 |
| Pentachlorophenol | Ave | 0.1432 | 0.1286 | 0.0500 | 18.0 | 20.0 | -10.2 | 20.0 |
| Phenanthrene | Ave | 1.177 | 1.137 | 0.7000 | 9.66 | 10.0 | -3.4 | 20.0 |
| Anthracene | Ave | 1.190 | 1.146 | 0.7000 | 9.63 | 10.0 | -3.7 | 20.0 |
| Carbazole | Ave | 1.127 | 1.044 | 0.0100 | 9.26 | 10.0 | -7.4 | 20.0 |
| Di-n-butyl phthalate | Ave | 1.380 | 1.388 | 0.0100 | 10.1 | 10.0 | 0.6 | 20.0 |
| Fluoranthene | Ave | 1.321 | 1.235 | 0.6000 | 9.35 | 10.0 | -6.5 | 20.0 |
| Benzidine | Ave | 0.5843 | 0.3860 | | 13.2 | 20.0 | -33.9* | 20.0 |
| Pyrene | Ave | 1.135 | 1.113 | 0.6000 | 9.81 | 10.0 | -1.9 | 20.0 |
| Butyl benzyl phthalate | Ave | 0.5142 | 0.5301 | 0.0100 | 10.3 | 10.0 | 3.1 | 20.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.6579 | 0.7290 | 0.0100 | 11.1 | 10.0 | 10.8 | 20.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3952 | 0.3826 | 0.0100 | 19.4 | 20.0 | -3.2 | 20.0 |
| Benzo[a]anthracene | Ave | 1.126 | 1.060 | 0.8000 | 9.41 | 10.0 | -5.9 | 20.0 |
| Chrysene | Ave | 1.076 | 0.9787 | 0.7000 | 9.10 | 10.0 | -9.0 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206256/2  Calibration Date: 11/11/2015 09:34

Instrument ID: A4HP10  Calib Start Date: 10/02/2015 10:48

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/02/2015 14:04

Lab File ID: 51111002.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Ave | 1.371 | 1.497 | 0.0100 | 10.9 | 10.0 | 9.2 | 20.0 |
| Benzo[b]fluoranthene | Ave | 1.128 | 1.167 | 0.7000 | 10.3 | 10.0 | 3.5 | 20.0 |
| Benzo[k]fluoranthene | Ave | 1.274 | 1.217 | 0.7000 | 9.55 | 10.0 | -4.5 | 20.0 |
| Benzo[a]pyrene | Ave | 1.120 | 1.130 | 0.7000 | 10.1 | 10.0 | 0.9 | 20.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.233 | 1.265 | 0.5000 | 10.3 | 10.0 | 2.6 | 20.0 |
| Dibenz(a,h)anthracene | Ave | 1.024 | 1.075 | 0.4000 | 10.5 | 10.0 | 5.0 | 20.0 |
| Benzo[g,h,i]perylene | Ave | 1.092 | 1.055 | 0.5000 | 9.66 | 10.0 | -3.4 | 20.0 |
| 2-Fluorophenol (Surr) | Ave | 1.486 | 1.313 | | 8.83 | 10.0 | -11.7 | 20.0 |
| Phenol-d5 (Surr) | Ave | 1.814 | 1.632 | | 9.00 | 10.0 | -10.0 | 20.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.3982 | 0.3544 | | 8.90 | 10.0 | -11.0 | 20.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.297 | 1.228 | | 9.47 | 10.0 | -5.3 | 20.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1891 | 0.1700 | | 8.99 | 10.0 | -10.1 | 20.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7619 | 0.7712 | | 10.1 | 10.0 | 1.2 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204569/11 Calibration Date: 10/30/2015 17:08

Instrument ID: A4HP9 Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS ID: 0.45(mm) Calib End Date: 10/30/2015 16:42

Lab File ID: 51030011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.7606 | 0.7555 | | 9.93 | 10.0 | -0.7 | 30.0 |
| N-Nitrosodimethylamine | Ave | 1.097 | 0.9653 | | 8.80 | 10.0 | -12.0 | 30.0 |
| Pyridine | Ave | 2.018 | 1.970 | | 9.76 | 10.0 | -2.4 | 30.0 |
| Benzaldehyde | Ave | 1.420 | 1.308 | 0.0100 | 18.4 | 20.0 | -7.9 | 30.0 |
| Phenol | Ave | 2.099 | 2.079 | 0.8000 | 9.91 | 10.0 | -0.9 | 30.0 |
| Aniline | Ave | 2.617 | 2.563 | | 9.79 | 10.0 | -2.1 | 30.0 |
| Bis(2-chloroethyl)ether | Ave | 1.803 | 1.702 | 0.7000 | 9.44 | 10.0 | -5.6 | 30.0 |
| 2-Chlorophenol | Ave | 1.517 | 1.526 | 0.8000 | 10.1 | 10.0 | 0.6 | 30.0 |
| n-Decane | Ave | 1.758 | 1.658 | | 9.44 | 10.0 | -5.6 | 30.0 |
| 1,3-Dichlorobenzene | Ave | 1.573 | 1.505 | | 9.57 | 10.0 | -4.3 | 30.0 |
| 1,4-Dichlorobenzene | Ave | 1.602 | 1.552 | | 9.69 | 10.0 | -3.1 | 30.0 |
| Benzyl alcohol | Ave | 1.029 | 1.068 | | 10.4 | 10.0 | 3.8 | 30.0 |
| 1,2-Dichlorobenzene | Ave | 1.540 | 1.462 | | 9.49 | 10.0 | -5.1 | 30.0 |
| 2-Methylphenol | Ave | 1.570 | 1.558 | 0.7000 | 9.93 | 10.0 | -0.7 | 30.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.024 | 1.968 | | 9.73 | 10.0 | -2.7 | 30.0 |
| Indene | Ave | 2.688 | 2.473 | | 18.4 | 20.0 | -8.0 | 30.0 |
| 3 & 4 Methylphenol | Ave | 1.620 | 1.612 | | 9.96 | 10.0 | -0.4 | 30.0 |
| N-Nitrosodi-n-propylamine | Ave | 1.139 | 1.142 | 0.5000 | 10.0 | 10.0 | 0.3 | 30.0 |
| Acetophenone | Ave | 2.353 | 2.256 | 0.0100 | 9.59 | 10.0 | -4.1 | 30.0 |
| Hexachloroethane | Ave | 0.6683 | 0.6557 | 0.3000 | 9.81 | 10.0 | -1.9 | 30.0 |
| Nitrobenzene | Ave | 0.4459 | 0.4395 | 0.2000 | 9.86 | 10.0 | -1.4 | 30.0 |
| Isophorone | Ave | 0.7626 | 0.7927 | 0.4000 | 10.4 | 10.0 | 3.9 | 30.0 |
| 2-Nitrophenol | Ave | 0.1660 | 0.1808 | 0.1000 | 10.9 | 10.0 | 8.9 | 30.0 |
| 2,4-Dimethylphenol | Ave | 0.3818 | 0.3725 | 0.2000 | 9.76 | 10.0 | -2.4 | 30.0 |
| Benzoic acid | Lin1 | | 0.2152 | | 20.9 | 20.0 | 4.4 | 30.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4539 | 0.4473 | 0.3000 | 9.86 | 10.0 | -1.4 | 30.0 |
| 2,4-Dichlorophenol | Ave | 0.2489 | 0.2521 | 0.2000 | 10.1 | 10.0 | 1.3 | 30.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.2911 | 0.2788 | | 9.58 | 10.0 | -4.2 | 30.0 |
| Naphthalene | Ave | 1.093 | 1.042 | 0.7000 | 9.53 | 10.0 | -4.7 | 30.0 |
| 4-Chloroaniline | Ave | 0.4651 | 0.4784 | 0.0100 | 10.3 | 10.0 | 2.9 | 30.0 |
| 2,6-Dichlorophenol | Ave | 0.2535 | 0.2551 | | 10.1 | 10.0 | 0.6 | 30.0 |
| Hexachlorobutadiene | Ave | 0.1548 | 0.1494 | 0.0100 | 9.65 | 10.0 | -3.5 | 30.0 |
| Caprolactam | Ave | 0.1118 | 0.1204 | 0.0100 | 21.5 | 20.0 | 7.7 | 30.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3177 | 0.3422 | 0.2000 | 10.8 | 10.0 | 7.7 | 30.0 |
| 2-Methylnaphthalene | Ave | 0.6768 | 0.6510 | 0.4000 | 9.62 | 10.0 | -3.8 | 30.0 |
| 1-Methylnaphthalene | Ave | 0.6004 | 0.6115 | | 10.2 | 10.0 | 1.9 | 30.0 |
| Hexachlorocyclopentadiene | Ave | 0.2585 | 0.2932 | 0.0500 | 11.3 | 10.0 | 13.4 | 30.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4935 | 0.4572 | 0.0100 | 9.26 | 10.0 | -7.4 | 30.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3377 | 0.3579 | 0.2000 | 10.6 | 10.0 | 6.0 | 30.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3511 | 0.3667 | 0.2000 | 10.4 | 10.0 | 4.4 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204569/11  Calibration Date: 10/30/2015 17:08

Instrument ID: A4HP9  Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS  ID: 0.45(mm)  Calib End Date: 10/30/2015 16:42

Lab File ID: 51030011.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.542 | 1.453 | 0.0100 | 9.42 | 10.0 | -5.8 | 30.0 |
| 2-Chloronaphthalene | Ave | 1.162 | 1.117 | 0.8000 | 9.62 | 10.0 | -3.8 | 30.0 |
| 2-Nitroaniline | Ave | 0.4031 | 0.4426 | 0.0100 | 11.0 | 10.0 | 9.8 | 30.0 |
| Dimethyl phthalate | Ave | 1.213 | 1.228 | 0.0100 | 10.1 | 10.0 | 1.2 | 30.0 |
| 1,3-Dinitrobenzene | Ave | 0.1896 | 0.2188 | | 11.5 | 10.0 | 15.4 | 30.0 |
| 2,6-Dinitrotoluene | Ave | 0.2814 | 0.2991 | | 10.6 | 10.0 | 6.3 | 30.0 |
| Acenaphthylene | Ave | 1.895 | 1.849 | 0.9000 | 9.76 | 10.0 | -2.4 | 30.0 |
| 3-Nitroaniline | Ave | 0.3457 | 0.3664 | 0.0100 | 10.6 | 10.0 | 6.0 | 30.0 |
| 2,4-Dinitrophenol | Lin | | 0.1277 | 0.0100 | 19.3 | 20.0 | -3.4 | 30.0 |
| Acenaphthene | Ave | 1.225 | 1.185 | 0.9000 | 9.67 | 10.0 | -3.3 | 30.0 |
| 4-Nitrophenol | Ave | 0.2054 | 0.2330 | | 22.7 | 20.0 | 13.4 | 30.0 |
| 2,4-Dinitrotoluene | Ave | 0.3558 | 0.3806 | 0.2000 | 10.7 | 10.0 | 7.0 | 30.0 |
| Dibenzofuran | Ave | 1.646 | 1.554 | 0.8000 | 9.44 | 10.0 | -5.6 | 30.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.2627 | 0.2874 | 0.0100 | 10.9 | 10.0 | 9.4 | 30.0 |
| Diethyl phthalate | Ave | 1.206 | 1.131 | 0.0100 | 9.38 | 10.0 | -6.2 | 30.0 |
| Hexadecane | Ave | 0.8758 | 0.8247 | | 9.42 | 10.0 | -5.8 | 30.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.5909 | 0.5890 | 0.4000 | 9.97 | 10.0 | -0.3 | 30.0 |
| 4-Nitroaniline | Ave | 0.3467 | 0.3816 | 0.0100 | 11.0 | 10.0 | 10.1 | 30.0 |
| Fluorene | Ave | 1.288 | 1.273 | 0.9000 | 9.89 | 10.0 | -1.1 | 30.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1284 | 0.0100 | 21.0 | 20.0 | 4.8 | 30.0 |
| N-Nitrosodiphenylamine | Ave | 0.5262 | 0.5266 | 0.0100 | 20.0 | 20.0 | 0.0 | 30.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 1.014 | 1.113 | | 11.0 | 10.0 | 9.8 | 30.0 |
| Azobenzene | Ave | 1.014 | 1.112 | | 11.0 | 10.0 | 9.7 | 30.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.1862 | 0.1987 | 0.1000 | 10.7 | 10.0 | 6.7 | 30.0 |
| Hexachlorobenzene | Ave | 0.1906 | 0.1906 | 0.1000 | 10.0 | 10.0 | 0.0 | 30.0 |
| Atrazine | Ave | 0.1757 | 0.1916 | 0.0100 | 21.8 | 20.0 | 9.0 | 30.0 |
| n-Octadecane | Ave | 0.5128 | 0.5469 | | 10.7 | 10.0 | 6.6 | 30.0 |
| Pentachlorophenol | Lin1 | | 0.1354 | 0.0500 | 22.2 | 20.0 | 11.1 | 30.0 |
| Phenanthrene | Ave | 1.179 | 1.105 | 0.7000 | 9.38 | 10.0 | -6.2 | 30.0 |
| Anthracene | Ave | 1.143 | 1.148 | 0.7000 | 10.0 | 10.0 | 0.4 | 30.0 |
| Carbazole | Ave | 0.996 | 1.069 | 0.0100 | 10.7 | 10.0 | 7.3 | 30.0 |
| Di-n-butyl phthalate | Ave | 1.181 | 1.389 | 0.0100 | 11.8 | 10.0 | 17.6 | 30.0 |
| Fluoranthene | Ave | 1.125 | 1.216 | 0.6000 | 10.8 | 10.0 | 8.0 | 30.0 |
| Benzidine | Lin1 | | 0.6541 | | 20.2 | 20.0 | 0.8 | 30.0 |
| Pyrene | Ave | 1.165 | 1.126 | 0.6000 | 9.66 | 10.0 | -3.4 | 30.0 |
| Butyl benzyl phthalate | Ave | 0.5273 | 0.5758 | 0.0100 | 10.9 | 10.0 | 9.2 | 30.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.7086 | 0.7773 | 0.0100 | 11.0 | 10.0 | 9.7 | 30.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3426 | 0.3569 | 0.0100 | 20.8 | 20.0 | 4.2 | 30.0 |
| Benzo[a]anthracene | Ave | 1.114 | 1.131 | 0.8000 | 10.2 | 10.0 | 1.6 | 30.0 |
| Chrysene | Ave | 1.127 | 1.120 | 0.7000 | 9.93 | 10.0 | -0.7 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204569/11 Calibration Date: 10/30/2015 17:08

Instrument ID: A4HP9 Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS ID: 0.45(mm) Calib End Date: 10/30/2015 16:42

Lab File ID: 51030011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Lin1 | | 1.598 | 0.0100 | 10.3 | 10.0 | 2.7 | 30.0 |
| Benzo[b]fluoranthene | Ave | 1.034 | 1.155 | 0.7000 | 11.2 | 10.0 | 11.8 | 30.0 |
| Benzo[k]fluoranthene | Ave | 1.334 | 1.317 | 0.7000 | 9.87 | 10.0 | -1.3 | 30.0 |
| Benzo[a]pyrene | Ave | 1.098 | 1.155 | 0.7000 | 10.5 | 10.0 | 5.2 | 30.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.106 | 1.175 | 0.5000 | 10.6 | 10.0 | 6.2 | 30.0 |
| Dibenz(a,h)anthracene | Ave | 0.8758 | 0.9878 | 0.4000 | 11.3 | 10.0 | 12.8 | 30.0 |
| Benzo[g,h,i]perylene | Ave | 0.9937 | 1.012 | 0.5000 | 10.2 | 10.0 | 1.9 | 30.0 |
| 2-Fluorophenol (Surr) | Ave | 1.509 | 1.588 | | 10.5 | 10.0 | 5.2 | 30.0 |
| Phenol-d5 (Surr) | Ave | 1.961 | 2.063 | | 10.5 | 10.0 | 5.2 | 30.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.4689 | 0.4903 | | 10.5 | 10.0 | 4.6 | 30.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.261 | 1.273 | | 10.1 | 10.0 | 1.0 | 30.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1401 | 0.1609 | | 11.5 | 10.0 | 14.8 | 30.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7313 | 0.7754 | | 10.6 | 10.0 | 6.0 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206233/2  Calibration Date: 11/11/2015 07:49

Instrument ID: A4HP9  Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS  ID: 0.45(mm)  Calib End Date: 10/30/2015 16:42

Lab File ID: 51111002.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| N-Nitrosodimethylamine | Ave | 1.097 | 1.401 | | 12.8 | 10.0 | 27.7* | 20.0 |
| Pyridine | Ave | 2.018 | 2.368 | | 11.7 | 10.0 | 17.4 | 20.0 |
| Benzaldehyde | Ave | 1.420 | 0.9711 | 0.0100 | 13.7 | 20.0 | -31.6* | 20.0 |
| Phenol | Ave | 2.099 | 2.304 | 0.8000 | 11.0 | 10.0 | 9.8 | 20.0 |
| Aniline | Ave | 2.617 | 2.827 | | 10.8 | 10.0 | 8.0 | 20.0 |
| Bis(2-chloroethyl)ether | Ave | 1.803 | 2.052 | 0.7000 | 11.4 | 10.0 | 13.8 | 20.0 |
| 2-Chlorophenol | Ave | 1.517 | 1.503 | 0.8000 | 9.91 | 10.0 | -0.9 | 20.0 |
| n-Decane | Ave | 1.758 | 2.188 | | 12.4 | 10.0 | 24.5* | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.573 | 1.567 | | 9.97 | 10.0 | -0.3 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.602 | 1.600 | | 9.99 | 10.0 | -0.1 | 20.0 |
| Benzyl alcohol | Ave | 1.029 | 1.117 | | 10.9 | 10.0 | 8.6 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.540 | 1.527 | | 9.91 | 10.0 | -0.9 | 20.0 |
| 2-Methylphenol | Ave | 1.570 | 1.556 | 0.7000 | 9.91 | 10.0 | -0.9 | 20.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.024 | 2.359 | | 11.7 | 10.0 | 16.6 | 20.0 |
| Indene | Ave | 2.688 | 2.657 | | 19.8 | 20.0 | -1.1 | 20.0 |
| 3 & 4 Methylphenol | Ave | 1.620 | 1.541 | | 9.51 | 10.0 | -4.9 | 20.0 |
| N-Nitrosodi-n-propylamine | Ave | 1.139 | 1.450 | 0.5000 | 12.7 | 10.0 | 27.4* | 20.0 |
| Acetophenone | Ave | 2.353 | 2.244 | 0.0100 | 9.54 | 10.0 | -4.6 | 20.0 |
| Hexachloroethane | Ave | 0.6683 | 0.6695 | 0.3000 | 10.0 | 10.0 | 0.2 | 20.0 |
| Nitrobenzene | Ave | 0.4459 | 0.4894 | 0.2000 | 11.0 | 10.0 | 9.8 | 20.0 |
| Isophorone | Ave | 0.7626 | 0.8151 | 0.4000 | 10.7 | 10.0 | 6.9 | 20.0 |
| 2-Nitrophenol | Ave | 0.1660 | 0.1725 | 0.1000 | 10.4 | 10.0 | 3.9 | 20.0 |
| 2,4-Dimethylphenol | Ave | 0.3818 | 0.3609 | 0.2000 | 9.45 | 10.0 | -5.5 | 20.0 |
| Benzoic acid | Lin1 | | 0.1419 | | 14.3 | 20.0 | -28.4* | 20.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4539 | 0.4711 | 0.3000 | 10.4 | 10.0 | 3.8 | 20.0 |
| 2,4-Dichlorophenol | Ave | 0.2489 | 0.2500 | 0.2000 | 10.0 | 10.0 | 0.4 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.2911 | 0.2488 | | 8.54 | 10.0 | -14.6 | 20.0 |
| Naphthalene | Ave | 1.093 | 0.9231 | 0.7000 | 8.44 | 10.0 | -15.6 | 20.0 |
| 4-Chloroaniline | Ave | 0.4651 | 0.3947 | 0.0100 | 8.49 | 10.0 | -15.1 | 20.0 |
| 2,6-Dichlorophenol | Ave | 0.2535 | 0.2456 | | 9.69 | 10.0 | -3.1 | 20.0 |
| Hexachlorobutadiene | Ave | 0.1548 | 0.1309 | 0.0100 | 8.46 | 10.0 | -15.4 | 20.0 |
| Caprolactam | Ave | 0.1118 | 0.1051 | 0.0100 | 18.8 | 20.0 | -6.0 | 20.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3177 | 0.3036 | 0.2000 | 9.56 | 10.0 | -4.4 | 20.0 |
| 2-Methylnaphthalene | Ave | 0.6768 | 0.6149 | 0.4000 | 9.09 | 10.0 | -9.1 | 20.0 |
| 1-Methylnaphthalene | Ave | 0.6004 | 0.5452 | | 9.08 | 10.0 | -9.2 | 20.0 |
| Hexachlorocyclopentadiene | Ave | 0.2585 | 0.2068 | 0.0500 | 8.00 | 10.0 | -20.0 | 20.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4935 | 0.4285 | 0.0100 | 8.68 | 10.0 | -13.2 | 20.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3377 | 0.3103 | 0.2000 | 9.19 | 10.0 | -8.1 | 20.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3511 | 0.3097 | 0.2000 | 8.82 | 10.0 | -11.8 | 20.0 |
| 1,1'-Biphenyl | Ave | 1.542 | 1.307 | 0.0100 | 8.48 | 10.0 | -15.2 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206233/2 | Calibration Date: 11/11/2015 07:49

Instrument ID: A4HP9 | Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS ID: 0.45(mm) | Calib End Date: 10/30/2015 16:42

Lab File ID: 51111002.D | Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 2-Chloronaphthalene | Ave | 1.162 | 1.068 | 0.8000 | 9.19 | 10.0 | -8.1 | 20.0 |
| 2-Nitroaniline | Ave | 0.4031 | 0.4647 | 0.0100 | 11.5 | 10.0 | 15.3 | 20.0 |
| Dimethyl phthalate | Ave | 1.213 | 1.112 | 0.0100 | 9.17 | 10.0 | -8.3 | 20.0 |
| 1,3-Dinitrobenzene | Ave | 0.1896 | 0.2036 | | 10.7 | 10.0 | 7.4 | 20.0 |
| 2,6-Dinitrotoluene | Ave | 0.2814 | 0.2741 | | 9.74 | 10.0 | -2.6 | 20.0 |
| Acenaphthylene | Ave | 1.895 | 1.621 | 0.9000 | 8.55 | 10.0 | -14.5 | 20.0 |
| 3-Nitroaniline | Ave | 0.3457 | 0.3215 | 0.0100 | 9.30 | 10.0 | -7.0 | 20.0 |
| 2,4-Dinitrophenol | Lin | | 0.1118 | 0.0100 | 17.4 | 20.0 | -13.2 | 20.0 |
| Acenaphthene | Ave | 1.225 | 1.006 | 0.9000 | 8.21 | 10.0 | -17.9 | 20.0 |
| 4-Nitrophenol | Ave | 0.2054 | 0.1907 | | 18.6 | 20.0 | -7.2 | 20.0 |
| 2,4-Dinitrotoluene | Ave | 0.3558 | 0.3600 | 0.2000 | 10.1 | 10.0 | 1.2 | 20.0 |
| Dibenzofuran | Ave | 1.646 | 1.411 | 0.8000 | 8.57 | 10.0 | -14.3 | 20.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.2627 | 0.2179 | 0.0100 | 8.29 | 10.0 | -17.1 | 20.0 |
| Diethyl phthalate | Ave | 1.206 | 0.9799 | 0.0100 | 8.12 | 10.0 | -18.8 | 20.0 |
| Hexadecane | Ave | 0.8758 | 0.8735 | | 9.97 | 10.0 | -0.3 | 20.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.5909 | 0.5060 | 0.4000 | 8.56 | 10.0 | -14.4 | 20.0 |
| 4-Nitroaniline | Ave | 0.3467 | 0.3204 | 0.0100 | 9.24 | 10.0 | -7.6 | 20.0 |
| Fluorene | Ave | 1.288 | 1.141 | 0.9000 | 8.86 | 10.0 | -11.4 | 20.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1238 | 0.0100 | 20.3 | 20.0 | 1.4 | 20.0 |
| N-Nitrosodiphenylamine | Ave | 0.5262 | 0.5308 | 0.0100 | 20.2 | 20.0 | 0.9 | 20.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 1.014 | 1.110 | | 10.9 | 10.0 | 9.5 | 20.0 |
| Azobenzene | Ave | 1.014 | 1.111 | | 11.0 | 10.0 | 9.5 | 20.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.1862 | 0.1790 | 0.1000 | 9.61 | 10.0 | -3.9 | 20.0 |
| Hexachlorobenzene | Ave | 0.1906 | 0.1928 | 0.1000 | 10.1 | 10.0 | 1.2 | 20.0 |
| Atrazine | Ave | 0.1757 | 0.1904 | 0.0100 | 21.7 | 20.0 | 8.3 | 20.0 |
| n-Octadecane | Ave | 0.5128 | 0.6714 | | 13.1 | 10.0 | 30.9* | 20.0 |
| Pentachlorophenol | Lin1 | | 0.1117 | 0.0500 | 18.5 | 20.0 | -7.7 | 20.0 |
| Phenanthrene | Ave | 1.179 | 1.084 | 0.7000 | 9.19 | 10.0 | -8.1 | 20.0 |
| Anthracene | Ave | 1.143 | 1.091 | 0.7000 | 9.54 | 10.0 | -4.6 | 20.0 |
| Carbazole | Ave | 0.996 | 1.073 | 0.0100 | 10.8 | 10.0 | 7.7 | 20.0 |
| Di-n-butyl phthalate | Ave | 1.181 | 1.258 | 0.0100 | 10.6 | 10.0 | 6.5 | 20.0 |
| Fluoranthene | Ave | 1.125 | 1.144 | 0.6000 | 10.2 | 10.0 | 1.7 | 20.0 |
| Benzidine | Lin1 | | 0.4427 | | 13.9 | 20.0 | -30.7* | 20.0 |
| Pyrene | Ave | 1.165 | 1.127 | 0.6000 | 9.67 | 10.0 | -3.3 | 20.0 |
| Butyl benzyl phthalate | Ave | 0.5273 | 0.5343 | 0.0100 | 10.1 | 10.0 | 1.3 | 20.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.7086 | 0.7187 | 0.0100 | 10.1 | 10.0 | 1.4 | 20.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3426 | 0.3520 | 0.0100 | 20.5 | 20.0 | 2.7 | 20.0 |
| Benzo[a]anthracene | Ave | 1.114 | 1.008 | 0.8000 | 9.05 | 10.0 | -9.5 | 20.0 |
| Chrysene | Ave | 1.127 | 0.9562 | 0.7000 | 8.48 | 10.0 | -15.2 | 20.0 |
| Di-n-octyl phthalate | Lin1 | | 1.464 | 0.0100 | 9.43 | 10.0 | -5.7 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206233/2  Calibration Date: 11/11/2015 07:49

Instrument ID: A4HP9  Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS  ID: 0.45(mm)  Calib End Date: 10/30/2015 16:42

Lab File ID: 51111002.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Benzo[b]fluoranthene | Ave | 1.034 | 1.122 | 0.7000 | 10.9 | 10.0 | 8.6 | 20.0 |
| Benzo[k]fluoranthene | Ave | 1.334 | 1.254 | 0.7000 | 9.40 | 10.0 | -6.0 | 20.0 |
| Benzo[a]pyrene | Ave | 1.098 | 1.113 | 0.7000 | 10.1 | 10.0 | 1.3 | 20.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.106 | 1.182 | 0.5000 | 10.7 | 10.0 | 6.8 | 20.0 |
| Dibenz(a,h)anthracene | Ave | 0.8758 | 0.9640 | 0.4000 | 11.0 | 10.0 | 10.1 | 20.0 |
| Benzo[g,h,i]perylene | Ave | 0.9937 | 0.9829 | 0.5000 | 9.89 | 10.0 | -1.1 | 20.0 |
| 1,4-Dioxane | Ave | 0.7606 | | | 0.000490 | 10.0 | | |
| 2-Fluorophenol (Surr) | Ave | 1.509 | 1.580 | | 10.5 | 10.0 | 4.7 | 20.0 |
| Phenol-d5 (Surr) | Ave | 1.961 | 2.103 | | 10.7 | 10.0 | 7.2 | 20.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.4689 | 0.4936 | | 10.5 | 10.0 | 5.3 | 20.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.261 | 1.114 | | 8.84 | 10.0 | -11.6 | 20.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1401 | 0.1104 | | 7.88 | 10.0 | -21.2* | 20.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7313 | 0.7199 | | 9.84 | 10.0 | -1.6 | 20.0 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205786/23-A

Matrix: Solid | Lab File ID: 51111004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:27

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 3.5 | U | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 9.5 | U | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 25 | U | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 8.9 | U | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 20 | U | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 20 | U | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 21 | U | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 17 | U | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 21 | U | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 0.45 | U | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 8.2 | U | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 0.50 | U | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 11 | U | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 9.1 | U | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 8.3 | U | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 18 | U | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 16 | U | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 9.2 | U | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 13 | U | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 21 | U | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 17 | U | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 13 | U | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 26 | U | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 17 | U | 330 | 17 |
| 83-32-9 | Acenaphthene | 0.76 | U | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 0.35 | U | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 9.2 | U | 100 | 9.2 |
| 120-12-7 | Anthracene | 0.78 | U | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 9.1 | U | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 12 | U | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 0.63 | U | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 0.64 | U | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 0.59 | U | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.35 | U | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205786/23-A

Matrix: Solid | Lab File ID: 51111004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:27

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.68 | U | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 22 | U | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.0 | U | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 19 | U | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 10 | U | 70 | 10 |
| 105-60-2 | Caprolactam | 37 | U | 330 | 37 |
| 86-74-8 | Carbazole | 27 | U | 50 | 27 |
| 218-01-9 | Chrysene | 1.1 | U | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.66 | U | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 0.66 | U | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 16 | U | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 17 | U | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 15 | U | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 7.9 | U | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 0.55 | U | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 0.53 | U | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 2.1 | U | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 5.6 | U | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 8.1 | U | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 9.0 | U | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.35 | U | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 13 | U | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 6.3 | U | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 21 | U | 50 | 21 |
| 91-20-3 | Naphthalene | 0.82 | U | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 2.2 | U | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 9.1 | U | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 0.73 | U | 6.7 | 0.73 |
| 108-95-2 | Phenol | 7.3 | U | 50 | 7.3 |
| 129-00-0 | Pyrene | 0.44 | U | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 20 | U | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205786/23-A

Matrix: Solid | Lab File ID: 51111004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:27

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 80 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 66 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 65 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 66 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 68 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 48 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205866/8-A

Matrix: Solid | Lab File ID: 51111006.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 09:54

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 3.5 | U | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 9.5 | U | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 25 | U | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 8.9 | U | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 20 | U | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 20 | U | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 21 | U | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 17 | U | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 21 | U | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 0.45 | U | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 8.2 | U | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 0.50 | U | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 11 | U | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 9.1 | U | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 8.3 | U | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 18 | U | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 16 | U | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 9.2 | U | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 13 | U | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 21 | U | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 17 | U | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 13 | U | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 26 | U | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 17 | U | 330 | 17 |
| 83-32-9 | Acenaphthene | 0.76 | U | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 0.35 | U | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 9.2 | U | 100 | 9.2 |
| 120-12-7 | Anthracene | 0.78 | U | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 9.1 | U | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 12 | U | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 0.63 | U | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 0.64 | U | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 0.59 | U | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.35 | U | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205866/8-A

Matrix: Solid | Lab File ID: 51111006.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 09:54

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.68 | U | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 22 | U | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.0 | U | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 19 | U | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 10 | U | 70 | 10 |
| 105-60-2 | Caprolactam | 37 | U | 330 | 37 |
| 86-74-8 | Carbazole | 27 | U | 50 | 27 |
| 218-01-9 | Chrysene | 1.1 | U | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.66 | U | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 0.66 | U | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 16 | U | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 17 | U | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 15 | U | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 7.9 | U | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 0.55 | U | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 0.53 | U | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 2.1 | U | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 5.6 | U | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 8.1 | U | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 9.0 | U | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.35 | U | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 13 | U | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 6.3 | U | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 21 | U | 50 | 21 |
| 91-20-3 | Naphthalene | 0.82 | U | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 2.2 | U | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 9.1 | U | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 0.73 | U | 6.7 | 0.73 |
| 108-95-2 | Phenol | 7.3 | U | 50 | 7.3 |
| 129-00-0 | Pyrene | 0.44 | U | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 20 | U | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205866/8-A

Matrix: Solid | Lab File ID: 51111006.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 09:54

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 87 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 88 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 78 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 81 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 76 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 49 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-205786/24-A

Matrix: Solid | Lab File ID: 51111005.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:51

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 525 | | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 611 | | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 428 | | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 319 | | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 532 | | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 462 | | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 481 | | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 605 | | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 563 | | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 523 | | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 528 | | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 550 | | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 576 | | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 596 | | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 584 | | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 900 | | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 497 | | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 716 | | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 571 | | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 571 | | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 395 | | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 549 | | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 588 | | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 1060 | | 330 | 17 |
| 83-32-9 | Acenaphthene | 533 | | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 532 | | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 551 | | 100 | 9.2 |
| 120-12-7 | Anthracene | 545 | | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 1240 | | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 1150 | | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 553 | | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 592 | | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 592 | | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 573 | | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-205786/24-A

Matrix: Solid | Lab File ID: 51111005.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:51

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 571 | | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 537 | | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 533 | | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 662 | | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 604 | | 70 | 10 |
| 105-60-2 | Caprolactam | 1110 | | 330 | 37 |
| 86-74-8 | Carbazole | 564 | | 50 | 27 |
| 218-01-9 | Chrysene | 550 | | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 620 | | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 548 | | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 618 | | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 572 | | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 583 | | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 628 | | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 539 | | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 550 | | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 553 | | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 559 | | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 240 | J | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 539 | | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 606 | | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 547 | | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 563 | | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 1180 | | 50 | 21 |
| 91-20-3 | Naphthalene | 537 | | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 550 | | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 584 | | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 547 | | 6.7 | 0.73 |
| 108-95-2 | Phenol | 542 | | 50 | 7.3 |
| 129-00-0 | Pyrene | 574 | | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 526 | | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-205786/24-A

Matrix: Solid | Lab File ID: 51111005.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:51

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 93 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 79 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 79 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 79 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 81 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 44 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-205866/9-A

Matrix: Solid | Lab File ID: 51111007.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:19

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 396 | | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 514 | | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 325 | | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 243 | | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 446 | | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 444 | | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 464 | | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 473 | | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 462 | | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 425 | | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 439 | | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 428 | | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 438 | | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 548 | | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 459 | | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 946 | | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 446 | | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 724 | | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 468 | | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 444 | | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 366 | | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 412 | | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 452 | | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 804 | | 330 | 17 |
| 83-32-9 | Acenaphthene | 386 | | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 400 | | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 424 | | 100 | 9.2 |
| 120-12-7 | Anthracene | 462 | | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 1140 | | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 1090 | | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 451 | | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 491 | | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 559 | | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 467 | | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-205866/9-A

Matrix: Solid | Lab File ID: 51111007.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:19

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 429 | | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 489 | | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 503 | | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 544 | | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 512 | | 70 | 10 |
| 105-60-2 | Caprolactam | 1000 | | 330 | 37 |
| 86-74-8 | Carbazole | 536 | | 50 | 27 |
| 218-01-9 | Chrysene | 442 | | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 513 | | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 406 | | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 418 | | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 455 | | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 538 | | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 448 | | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 500 | | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 426 | | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 496 | | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 392 | | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 244 | J | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 423 | | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 505 | | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 513 | | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 592 | | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 987 | | 50 | 21 |
| 91-20-3 | Naphthalene | 387 | | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 506 | | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 573 | | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 439 | | 6.7 | 0.73 |
| 108-95-2 | Phenol | 537 | | 50 | 7.3 |
| 129-00-0 | Pyrene | 481 | | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 432 | | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-205866/9-A

Matrix: Solid | Lab File ID: 51111007.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:19

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 76 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 75 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 76 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 72 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 64 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 33 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: 51111024.D

Analysis Method: 8270D | Date Collected: 11/05/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30.05(g) | Date Analyzed: 11/11/2015 16:52

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 511 | | 58 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 646 | | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 290 | | 180 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 318 | | 180 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 444 | | 180 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 275 | | 180 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 940 | | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 616 | | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 597 | | 230 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 546 | | 58 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 504 | | 58 | 9.6 |
| 91-57-6 | 2-Methylnaphthalene | 528 | | 7.8 | 0.58 |
| 95-48-7 | 2-Methylphenol | 391 | | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 709 | | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 593 | | 58 | 9.7 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 1180 | | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 533 | | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 1080 | | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 592 | | 58 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 231 | | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 391 | | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 537 | | 58 | 15 |
| 100-01-6 | 4-Nitroaniline | 570 | | 230 | 30 |
| 100-02-7 | 4-Nitrophenol | 383 | J | 390 | 20 |
| 83-32-9 | Acenaphthene | 501 | | 7.8 | 0.89 |
| 208-96-8 | Acenaphthylene | 518 | | 7.8 | 0.41 |
| 98-86-2 | Acetophenone | 542 | | 120 | 11 |
| 120-12-7 | Anthracene | 585 | | 7.8 | 0.91 |
| 1912-24-9 | Atrazine | 1450 | | 230 | 11 |
| 100-52-7 | Benzaldehyde | 1350 | | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 556 | | 7.8 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 623 | | 7.8 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 673 | | 7.8 | 0.69 |
| 191-24-2 | Benzo[g,h,i]perylene | 598 | | 7.8 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: 51111024.D

Analysis Method: 8270D | Date Collected: 11/05/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30.05(g) | Date Analyzed: 11/11/2015 16:52

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 562 | | 7.8 | 0.79 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 611 | | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 607 | | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 679 | | 82 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 649 | | 82 | 12 |
| 105-60-2 | Caprolactam | 1260 | | 390 | 43 |
| 86-74-8 | Carbazole | 666 | | 58 | 32 |
| 218-01-9 | Chrysene | 544 | | 7.8 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 637 | | 7.8 | 0.77 |
| 132-64-9 | Dibenzofuran | 536 | | 58 | 0.77 |
| 84-66-2 | Diethyl phthalate | 546 | | 82 | 19 |
| 131-11-3 | Dimethyl phthalate | 595 | | 82 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 681 | | 82 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 559 | | 82 | 9.2 |
| 206-44-0 | Fluoranthene | 621 | | 7.8 | 0.64 |
| 86-73-7 | Fluorene | 548 | | 7.8 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 623 | | 7.8 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 472 | | 58 | 6.5 |
| 77-47-4 | Hexachlorocyclopentadiene | 219 | J | 390 | 9.5 |
| 67-72-1 | Hexachloroethane | 526 | | 58 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 622 | | 7.8 | 0.41 |
| 78-59-1 | Isophorone | 651 | | 58 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 757 | | 58 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 1270 | | 58 | 25 |
| 91-20-3 | Naphthalene | 473 | | 7.8 | 0.96 |
| 98-95-3 | Nitrobenzene | 633 | | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 423 | | 180 | 11 |
| 85-01-8 | Phenanthrene | 557 | | 7.8 | 0.85 |
| 108-95-2 | Phenol | 395 | | 58 | 8.5 |
| 129-00-0 | Pyrene | 603 | | 7.8 | 0.51 |
| 15831-10-4 | 3 & 4 Methylphenol | 368 | J | 470 | 23 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: 51111024.D

Analysis Method: 8270D | Date Collected: 11/05/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30.05(g) | Date Analyzed: 11/11/2015 16:52

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 80 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 62 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 78 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 69 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 71 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 30 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD | Lab Sample ID: 240-57597-15 MSD

Matrix: Solid | Lab File ID: 51111025.D

Analysis Method: 8270D | Date Collected: 11/05/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.93(g) | Date Analyzed: 11/11/2015 17:17

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 419 | | 59 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 536 | | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 208 | | 180 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 307 | | 180 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 368 | | 180 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 286 | | 180 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 819 | | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 589 | | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 546 | | 230 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 455 | | 59 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 422 | | 59 | 9.6 |
| 91-57-6 | 2-Methylnaphthalene | 437 | | 7.8 | 0.59 |
| 95-48-7 | 2-Methylphenol | 416 | | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 646 | | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 492 | | 59 | 9.7 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 1210 | | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 551 | | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 996 | | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 532 | | 59 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 272 | | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 392 | | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 476 | | 59 | 15 |
| 100-01-6 | 4-Nitroaniline | 541 | | 230 | 30 |
| 100-02-7 | 4-Nitrophenol | 446 | | 390 | 20 |
| 83-32-9 | Acenaphthene | 432 | | 7.8 | 0.89 |
| 208-96-8 | Acenaphthylene | 443 | | 7.8 | 0.41 |
| 98-86-2 | Acetophenone | 442 | | 120 | 11 |
| 120-12-7 | Anthracene | 523 | | 7.8 | 0.91 |
| 1912-24-9 | Atrazine | 1300 | | 230 | 11 |
| 100-52-7 | Benzaldehyde | 1100 | | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 536 | | 7.8 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 563 | | 7.8 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 638 | | 7.8 | 0.69 |
| 191-24-2 | Benzo[g,h,i]perylene | 545 | | 7.8 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD | Lab Sample ID: 240-57597-15 MSD

Matrix: Solid | Lab File ID: 51111025.D

Analysis Method: 8270D | Date Collected: 11/05/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.93(g) | Date Analyzed: 11/11/2015 17:17

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 486 | | 7.8 | 0.80 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 497 | | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 528 | | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 661 | | 82 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 614 | | 82 | 12 |
| 105-60-2 | Caprolactam | 1170 | | 390 | 43 |
| 86-74-8 | Carbazole | 604 | | 59 | 32 |
| 218-01-9 | Chrysene | 521 | | 7.8 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 596 | | 7.8 | 0.77 |
| 132-64-9 | Dibenzofuran | 463 | | 59 | 0.77 |
| 84-66-2 | Diethyl phthalate | 507 | | 82 | 19 |
| 131-11-3 | Dimethyl phthalate | 543 | | 82 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 631 | | 82 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 517 | | 82 | 9.3 |
| 206-44-0 | Fluoranthene | 577 | | 7.8 | 0.64 |
| 86-73-7 | Fluorene | 486 | | 7.8 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 557 | | 7.8 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 395 | | 59 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 186 | J | 390 | 9.5 |
| 67-72-1 | Hexachloroethane | 411 | | 59 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 578 | | 7.8 | 0.41 |
| 78-59-1 | Isophorone | 540 | | 59 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 615 | | 59 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 1130 | | 59 | 25 |
| 91-20-3 | Naphthalene | 397 | | 7.8 | 0.96 |
| 98-95-3 | Nitrobenzene | 515 | | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 617 | | 180 | 11 |
| 85-01-8 | Phenanthrene | 510 | | 7.8 | 0.86 |
| 108-95-2 | Phenol | 418 | | 59 | 8.6 |
| 129-00-0 | Pyrene | 575 | | 7.8 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 392 | J | 470 | 23 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD | Lab Sample ID: 240-57597-15 MSD

Matrix: Solid | Lab File ID: 51111025.D

Analysis Method: 8270D | Date Collected: 11/05/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 29.93(g) | Date Analyzed: 11/11/2015 17:17

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 77 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 53 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 65 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 61 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 60 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 36 | | 10-110 |

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP10 Start Date: 10/02/2015 10:22

Analysis Batch Number: 200091 End Date: 10/02/2015 21:25

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| DFTPP 240-200091/1 | | 10/02/2015 10:22 | 1 | 51002001.D | DB-624 0.18(mm) |
| STD5 240-200091/2 IC | | 10/02/2015 10:48 | 1 | 51002002.D | DB-624 0.18(mm) |
| STD4 240-200091/3 IC | | 10/02/2015 11:12 | 1 | 51002003.D | DB-624 0.18(mm) |
| STD3 240-200091/4 IC | | 10/02/2015 11:37 | 1 | 51002004.D | DB-624 0.18(mm) |
| STD2 240-200091/5 IC | | 10/02/2015 12:01 | 1 | 51002005.D | DB-624 0.18(mm) |
| STD1 240-200091/6 IC | | 10/02/2015 12:25 | 1 | 51002006.D | DB-624 0.18(mm) |
| STD9 240-200091/7 IC | | 10/02/2015 12:50 | 1 | 51002007.D | DB-624 0.18(mm) |
| STD8 240-200091/8 IC | | 10/02/2015 13:15 | 1 | 51002008.D | DB-624 0.18(mm) |
| STD7 240-200091/9 IC | | 10/02/2015 13:39 | 1 | 51002009.D | DB-624 0.18(mm) |
| STD6 240-200091/10 ICIS | | 10/02/2015 14:04 | 1 | 51002010.D | DB-624 0.18(mm) |
| ICV 240-200091/11 | | 10/02/2015 14:28 | 1 | 51002011.D | DB-624 0.18(mm) |
| STD9 240-200091/12 IC | | 10/02/2015 14:53 | 1 | | DB-624 0.18(mm) |
| STD8 240-200091/13 IC | | 10/02/2015 15:17 | 1 | | DB-624 0.18(mm) |
| STD7 240-200091/14 IC | | 10/02/2015 15:42 | 1 | | DB-624 0.18(mm) |
| STD6 240-200091/15 IC | | 10/02/2015 16:06 | 1 | | DB-624 0.18(mm) |
| STD5 240-200091/16 IC | | 10/02/2015 16:31 | 1 | | DB-624 0.18(mm) |
| STD4 240-200091/17 IC | | 10/02/2015 16:55 | 1 | | DB-624 0.18(mm) |
| STD3 240-200091/18 IC | | 10/02/2015 17:20 | 1 | | DB-624 0.18(mm) |
| STD2 240-200091/19 IC | | 10/02/2015 17:45 | 1 | | DB-624 0.18(mm) |
| ICV 240-200091/20 | | 10/02/2015 18:09 | 1 | | DB-624 0.18(mm) |
| STD9 240-200091/21 IC | | 10/02/2015 18:33 | 1 | | DB-624 0.18(mm) |
| STD8 240-200091/22 IC | | 10/02/2015 18:58 | 1 | | DB-624 0.18(mm) |
| STD7 240-200091/23 IC | | 10/02/2015 19:22 | 1 | | DB-624 0.18(mm) |
| STD6 240-200091/24 IC | | 10/02/2015 19:47 | 1 | | DB-624 0.18(mm) |
| STD5 240-200091/25 IC | | 10/02/2015 20:12 | 1 | | DB-624 0.18(mm) |
| STD4 240-200091/26 IC | | 10/02/2015 20:36 | 1 | | DB-624 0.18(mm) |
| STD3 240-200091/27 IC | | 10/02/2015 21:00 | 1 | | DB-624 0.18(mm) |
| STD2 240-200091/28 IC | | 10/02/2015 21:25 | 1 | | DB-624 0.18(mm) |

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP10 Start Date: 11/11/2015 09:12

Analysis Batch Number: 206256 End Date: 11/11/2015 16:11

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| DFTPP 240-206256/1 | | 11/11/2015 09:12 | 1 | 51111001.D | DB-624 0.18(mm) |
| CCV 240-206256/2 CCVIS | | 11/11/2015 09:34 | 1 | 51111002.D | DB-624 0.18(mm) |
| LODV 240-206256/3 | | 11/11/2015 10:02 | 1 | | DB-624 0.18(mm) |
| MB 240-205786/23-A | | 11/11/2015 10:27 | 1 | 51111004.D | DB-624 0.18(mm) |
| LCS 240-205786/24-A | | 11/11/2015 10:51 | 1 | 51111005.D | DB-624 0.18(mm) |
| 240-57597-1 | F-VB-UST-1-14 | 11/11/2015 11:16 | 1 | 51111006.D | DB-624 0.18(mm) |
| 240-57597-2 | F-VB-UST-2-14 | 11/11/2015 11:40 | 1 | 51111007.D | DB-624 0.18(mm) |
| 240-57597-3 | F-VB-UST-3-14 | 11/11/2015 12:05 | 1 | 51111008.D | DB-624 0.18(mm) |
| 240-57597-4 | F-VB-UST-4-14 | 11/11/2015 12:30 | 1 | 51111009.D | DB-624 0.18(mm) |
| 240-57597-5 | F-VB-UST-5-14 | 11/11/2015 12:54 | 1 | 51111010.D | DB-624 0.18(mm) |
| 240-57597-6 | F-VB-UST-6-14 | 11/11/2015 13:19 | 1 | 51111011.D | DB-624 0.18(mm) |
| 240-57597-7 | F-VB-UST-7-14 | 11/11/2015 13:44 | 1 | 51111012.D | DB-624 0.18(mm) |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 14:08 | 1 | 51111013.D | DB-624 0.18(mm) |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 14:33 | 1 | 51111014.D | DB-624 0.18(mm) |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 14:57 | 1 | 51111015.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/11/2015 15:22 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/11/2015 15:47 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/11/2015 16:11 | 1 | | DB-624 0.18(mm) |

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP9 Start Date: 10/30/2015 12:57

Analysis Batch Number: 204569 End Date: 10/30/2015 17:08

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| DFTPP 240-204569/1 | | 10/30/2015 12:57 | 1 | 51030001.D | RXI-5SILMS 0.45(mm) |
| STD5 240-204569/2 IC | | 10/30/2015 13:19 | 1 | 51030002.D | RXI-5SILMS 0.45(mm) |
| STD4 240-204569/3 IC | | 10/30/2015 13:44 | 1 | 51030003.D | RXI-5SILMS 0.45(mm) |
| STD3 240-204569/4 IC | | 10/30/2015 14:10 | 1 | 51030004.D | RXI-5SILMS 0.45(mm) |
| STD2 240-204569/5 IC | | 10/30/2015 14:35 | 1 | 51030005.D | RXI-5SILMS 0.45(mm) |
| STD1 240-204569/6 IC | | 10/30/2015 15:01 | 1 | 51030006.D | RXI-5SILMS 0.45(mm) |
| STD9 240-204569/7 IC | | 10/30/2015 15:27 | 1 | 51030007.D | RXI-5SILMS 0.45(mm) |
| STD8 240-204569/8 IC | | 10/30/2015 15:52 | 1 | 51030008.D | RXI-5SILMS 0.45(mm) |
| STD7 240-204569/9 IC | | 10/30/2015 16:17 | 1 | 51030009.D | RXI-5SILMS 0.45(mm) |
| STD6 240-204569/10 ICIS | | 10/30/2015 16:42 | 1 | 51030010.D | RXI-5SILMS 0.45(mm) |
| ICV 240-204569/11 | | 10/30/2015 17:08 | 1 | 51030011.D | RXI-5SILMS 0.45(mm) |

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A4HP9 Start Date: 11/11/2015 07:27

Analysis Batch Number: 206233 End Date: 11/11/2015 17:17

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| DFTPP 240-206233/1 | | 11/11/2015 07:27 | 1 | 51111001.D | RXI-5SILMS 0.45(mm) |
| CCV 240-206233/2 CCVIS | | 11/11/2015 07:49 | 1 | 51111002.D | RXI-5SILMS 0.45(mm) |
| LODV 240-206233/3 | | 11/11/2015 08:16 | 1 | | RXI-5SILMS 0.45(mm) |
| LODV 240-206233/18 | | 11/11/2015 08:41 | 1 | | RXI-5SILMS 0.45(mm) |
| MB 240-205866/8-A | | 11/11/2015 09:54 | 1 | 51111006.D | RXI-5SILMS 0.45(mm) |
| LCS 240-205866/9-A | | 11/11/2015 10:19 | 1 | 51111007.D | RXI-5SILMS 0.45(mm) |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 14:50 | 1 | 51111019.D | RXI-5SILMS 0.45(mm) |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 15:14 | 1 | 51111020.D | RXI-5SILMS 0.45(mm) |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 15:39 | 1 | 51111021.D | RXI-5SILMS 0.45(mm) |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 16:03 | 1 | 51111022.D | RXI-5SILMS 0.45(mm) |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 16:28 | 1 | 51111023.D | RXI-5SILMS 0.45(mm) |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 16:52 | 1 | 51111024.D | RXI-5SILMS 0.45(mm) |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 17:17 | 1 | 51111025.D | RXI-5SILMS 0.45(mm) |

# GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205786  Batch Start Date: 11/09/15 10:35  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/10/15 02:35

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exBENZALDEHYD 00031 | exBNASPIKE 00053 | exBNASURR W 00028 | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-E-1 | F-VB-UST-1-14 | 3540C, 8270D | T | 29.99 g | 2 mL | | | 1 mL | |
| 240-57597-D-2 | F-VB-UST-2-14 | 3540C, 8270D | T | 30.16 g | 2 mL | | | 1 mL | |
| 240-57597-E-3 | F-VB-UST-3-14 | 3540C, 8270D | T | 30.03 g | 2 mL | | | 1 mL | |
| 240-57597-E-4 | F-VB-UST-4-14 | 3540C, 8270D | T | 29.84 g | 2 mL | | | 1 mL | |
| 240-57597-E-5 | F-VB-UST-5-14 | 3540C, 8270D | T | 30.28 g | 2 mL | | | 1 mL | |
| 240-57597-E-6 | F-VB-UST-6-14 | 3540C, 8270D | T | 30.15 g | 2 mL | | | 1 mL | |
| 240-57597-E-7 | F-VB-UST-7-14 | 3540C, 8270D | T | 29.53 g | 2 mL | | | 1 mL | |
| 240-57597-D-8 | F-VB-UST-8-14 | 3540C, 8270D | T | 30.20 g | 2 mL | | | 1 mL | |
| 240-57597-E-9 | F-VB-UST-9-14 | 3540C, 8270D | T | 29.67 g | 2 mL | | | 1 mL | |
| 240-57597-E-10 | F-VB-UST-10-14 | 3540C, 8270D | T | 29.78 g | 2 mL | | | 1 mL | |
| MB 240-205786/23 | | 3540C, 8270D | | 30 g | 2 mL | | | 1 mL | |
| LCS 240-205786/24 | | 3540C, 8270D | | 30 g | 2 mL | 1 mL | 1 mL | 1 mL | |

| Batch Notes | |
|---|---|
| Balance ID | 1339912 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2258961 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205866  Batch Start Date: 11/09/15 01:00  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/10/15 05:00

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exBENZALDEHYD 00031 | exBNASPIKE 00053 | exBNASURR W 00028 | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-E-11 | F-VB-UST-11-14 | 3540C, 8270D | T | 29.86 g | 2 mL | | | 1 mL | |
| 240-57597-E-12 | F-VB-UST-12-14 | 3540C, 8270D | T | 29.84 g | 2 mL | | | 1 mL | |
| 240-57597-E-13 | F-VB-UST-13-14 | 3540C, 8270D | T | 30.37 g | 2 mL | | | 1 mL | |
| 240-57597-D-14 | F-VB-UST-14-14 | 3540C, 8270D | T | 29.64 g | 2 mL | | | 1 mL | |
| 240-57597-D-15 | F-VB-UST-DUP | 3540C, 8270D | T | 30.09 g | 2 mL | | | 1 mL | |
| 240-57597-D-15 MS | F-VB-UST-DUP | 3540C, 8270D | T | 30.05 g | 2 mL | 1 mL | 1 mL | 1 mL | |
| 240-57597-D-15 MSD | F-VB-UST-DUP | 3540C, 8270D | T | 29.93 g | 2 mL | 1 mL | 1 mL | 1 mL | |
| MB 240-205866/8 | | 3540C, 8270D | | 30 g | 2 mL | | | 1 mL | |
| LCS 240-205866/9 | | 3540C, 8270D | | 30 g | 2 mL | 1 mL | 1 mL | 1 mL | |

| Batch Notes | |
|---|---|
| Balance ID | 1339907 |
| Na2SO4 Lot Number | 2204509 |
| Prep Solvent Lot # | 2258960 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8015C - GRO

## Nonhalogenated Organics using GC/FID - Modified (Gasoline Range Organics) by Method 8015C

FORM II
GC VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): RTX-VRX ID: 0.45(mm)

| Client Sample ID | Lab Sample ID | TFT1 # |
|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 78 |
| F-VB-UST-2-14 | 240-57597-2 | 79 |
| F-VB-UST-3-14 | 240-57597-3 | 77 |
| F-VB-UST-4-14 | 240-57597-4 | 71 |
| F-VB-UST-5-14 | 240-57597-5 | 71 |
| F-VB-UST-6-14 | 240-57597-6 | 76 |
| F-VB-UST-7-14 | 240-57597-7 | 78 |
| F-VB-UST-8-14 | 240-57597-8 | 73 |
| F-VB-UST-9-14 | 240-57597-9 | 78 |
| F-VB-UST-10-14 | 240-57597-10 | 79 |
| F-VB-UST-11-14 | 240-57597-11 | 79 |
| F-VB-UST-12-14 | 240-57597-12 | 78 |
| F-VB-UST-13-14 | 240-57597-13 | 75 |
| F-VB-UST-14-14 | 240-57597-14 | 80 |
| F-VB-UST-DUP | 240-57597-15 | 74 |
| | MB 240-206227/1-A | 73 |
| | MB 240-206444/1-A | 72 |
| | LCS 240-206227/2-A | 95 |
| | LCS 240-206444/2-A | 100 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 99 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 40 |

| | QC LIMITS |
|---|---|
| TFT = Trifluorotoluene (Surr) | 40-139 |

# Column to be used to flag recovery values

FORM II 8015C

FORM III
GC VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111108.D

Lab ID: LCS 240-206227/2-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 786 | 98 | 74-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111206.D

Lab ID: LCS 240-206444/2-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 787 | 98 | 74-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111126.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 928 | 54 U | 856 | 92 | 10-136 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111127.D

Lab ID: 240-57597-15 MSD Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 926 | 206 | 22 | 122 | 40 | 10-136 | F2 |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM IV
GC VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: MB 240-206227/1-A

Matrix: Solid  Date Extracted: 11/11/2015 06:58

Lab File ID:(1) A0111107.D  Lab File ID:(2)

Date Analyzed:(1) 11/11/2015 09:56  Date Analyzed:(2)

Instrument ID:(1) AFID  Instrument ID:(2)

GC Column:(1) RTX-VRX  ID: 0.45(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-206227/2-A | 11/11/2015 10:31 | |
| F-VB-UST-1-14 | 240-57597-1 | 11/11/2015 11:08 | |
| F-VB-UST-2-14 | 240-57597-2 | 11/11/2015 11:45 | |
| F-VB-UST-3-14 | 240-57597-3 | 11/11/2015 12:23 | |
| F-VB-UST-4-14 | 240-57597-4 | 11/11/2015 13:00 | |
| F-VB-UST-6-14 | 240-57597-6 | 11/11/2015 14:13 | |
| F-VB-UST-7-14 | 240-57597-7 | 11/11/2015 14:50 | |
| F-VB-UST-8-14 | 240-57597-8 | 11/11/2015 15:27 | |
| F-VB-UST-9-14 | 240-57597-9 | 11/11/2015 17:19 | |
| F-VB-UST-10-14 | 240-57597-10 | 11/11/2015 17:57 | |
| F-VB-UST-11-14 | 240-57597-11 | 11/11/2015 18:34 | |
| F-VB-UST-12-14 | 240-57597-12 | 11/11/2015 19:12 | |
| F-VB-UST-13-14 | 240-57597-13 | 11/11/2015 19:48 | |
| F-VB-UST-14-14 | 240-57597-14 | 11/11/2015 20:25 | |
| F-VB-UST-DUP | 240-57597-15 | 11/11/2015 21:03 | |
| F-VB-UST-DUP MS | 240-57597-15 MS | 11/11/2015 21:40 | |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 11/11/2015 22:17 | |

FORM IV
GC VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: MB 240-206444/1-A

Matrix: Solid  Date Extracted: 11/12/2015 05:59

Lab File ID:(1) A0111205.D  Lab File ID:(2)

Date Analyzed:(1) 11/12/2015 08:03  Date Analyzed:(2)

Instrument ID:(1) AFID  Instrument ID:(2)

GC Column:(1) RTX-VRX  ID: 0.45(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-206444/2-A | 11/12/2015 08:40 | |
| F-VB-UST-5-14 | 240-57597-5 | 11/12/2015 09:17 | |

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: STD4 240-194508/10  Date Analyzed: 08/24/2015 15:53

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): A0082110.D  Heated Purge: (Y/N) N

Calibration ID: 30361

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| INITIAL CALIBRATION SURROGATE | | | | 10.34 | | |
| UPPER LIMIT | | | | 10.44 | | |
| LOWER LIMIT | | | | 10.24 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| STD4 240-194508/10 ICRT | | 08/24/2015 15:53 | A0082110.D | 10.34 | | |
| ICV 240-194508/13 | | 08/24/2015 17:45 | A0082113.D | 10.34 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVRT 240-206211/4  Date Analyzed: 11/11/2015 08:08

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): A0111104.D  Heated Purge: (Y/N) N

Calibration ID: 30361

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 10.36 | | |
| UPPER LIMIT | | | | 10.46 | | |
| LOWER LIMIT | | | | 10.26 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCVRT 240-206211/4 | | 11/11/2015 08:08 | A0111104.D | 10.36 | | |
| CCV 240-206211/6 | | 11/11/2015 09:20 | A0111106.D | 10.37 | | |
| MB 240-206227/1-A | | 11/11/2015 09:56 | A0111107.D | 10.36 | | |
| LCS 240-206227/2-A | | 11/11/2015 10:31 | A0111108.D | 10.36 | | |
| 240-57597-1 | F-VB-UST-1-14 | 11/11/2015 11:08 | A0111109.D | 10.37 | | |
| 240-57597-2 | F-VB-UST-2-14 | 11/11/2015 11:45 | A0111110.D | 10.37 | | |
| 240-57597-3 | F-VB-UST-3-14 | 11/11/2015 12:23 | A0111111.D | 10.37 | | |
| 240-57597-4 | F-VB-UST-4-14 | 11/11/2015 13:00 | A0111112.D | 10.37 | | |
| 240-57597-6 | F-VB-UST-6-14 | 11/11/2015 14:13 | A0111114.D | 10.37 | | |
| 240-57597-7 | F-VB-UST-7-14 | 11/11/2015 14:50 | A0111115.D | 10.37 | | |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 15:27 | A0111116.D | 10.37 | | |
| CCV 240-206211/18 | | 11/11/2015 16:42 | A0111118.D | 10.37 | | |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 17:19 | A0111119.D | 10.37 | | |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 17:57 | A0111120.D | 10.37 | | |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 18:34 | A0111121.D | 10.37 | | |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 19:12 | A0111122.D | 10.37 | | |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 19:48 | A0111123.D | 10.37 | | |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 20:25 | A0111124.D | 10.37 | | |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 21:03 | A0111125.D | 10.37 | | |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 21:40 | A0111126.D | 10.37 | | |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 22:17 | A0111127.D | 10.37 | | |
| CCV 240-206211/28 | | 11/11/2015 22:54 | A0111128.D | 10.37 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVRT 240-206404/4  Date Analyzed: 11/12/2015 07:26

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): A0111204.D  Heated Purge: (Y/N) N

Calibration ID: 30361

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 10.37 | | |
| UPPER LIMIT | | | | 10.47 | | |
| LOWER LIMIT | | | | 10.27 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCVRT 240-206404/4 | | 11/12/2015 07:26 | A0111204.D | 10.37 | | |
| MB 240-206444/1-A | | 11/12/2015 08:03 | A0111205.D | 10.37 | | |
| LCS 240-206444/2-A | | 11/12/2015 08:40 | A0111206.D | 10.37 | | |
| 240-57597-5 | F-VB-UST-5-14 | 11/12/2015 09:17 | A0111207.D | 10.37 | | |
| CCV 240-206404/8 | | 11/12/2015 09:54 | A0111208.D | 10.37 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: A0111109.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 4.98(g) | Date Analyzed: 11/11/2015 11:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: A0111110.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 4.98(g) | Date Analyzed: 11/11/2015 11:45

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 79 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: A0111111.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 4.99(g) | Date Analyzed: 11/11/2015 12:23

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 56 | U | 120 | 56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 77 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14  Lab Sample ID: 240-57597-4

Matrix: Solid  Lab File ID: A0111112.D

Analysis Method: 8015C  Date Collected: 11/05/2015 09:06

Sample wt/vol: 4.99(g)  Date Analyzed: 11/11/2015 13:00

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX  ID: 0.45(mm)

% Moisture: 15.6  Level: (low/med) Low

Analysis Batch No.: 206211  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 71 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14  Lab Sample ID: 240-57597-5

Matrix: Solid  Lab File ID: A0111207.D

Analysis Method: 8015C  Date Collected: 11/05/2015 09:08

Sample wt/vol: 4.98(g)  Date Analyzed: 11/12/2015 09:17

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX  ID: 0.45(mm)

% Moisture: 15.2  Level: (low/med) Low

Analysis Batch No.: 206404  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 77 | J | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 71 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: A0111114.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.01(g) | Date Analyzed: 11/11/2015 14:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 76 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: A0111115.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 4.97(g) | Date Analyzed: 11/11/2015 14:50

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14  Lab Sample ID: 240-57597-8

Matrix: Solid  Lab File ID: A0111116.D

Analysis Method: 8015C  Date Collected: 11/05/2015 09:14

Sample wt/vol: 4.98(g)  Date Analyzed: 11/11/2015 15:27

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX  ID: 0.45(mm)

% Moisture: 18.0  Level: (low/med) Low

Analysis Batch No.: 206211  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 56 | U | 120 | 56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 73 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: A0111119.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.00(g) | Date Analyzed: 11/11/2015 17:19

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 57 | U | 120 | 57 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: A0111120.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.02(g) | Date Analyzed: 11/11/2015 17:57

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 53 | U | 120 | 53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 79 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: A0111121.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/11/2015 18:34

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 79 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: A0111122.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.04(g) | Date Analyzed: 11/11/2015 19:12

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: A0111123.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 4.99(g) | Date Analyzed: 11/11/2015 19:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 75 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: A0111124.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.02(g) | Date Analyzed: 11/11/2015 20:25

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 80 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: A0111125.D

Analysis Method: 8015C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.00(g) | Date Analyzed: 11/11/2015 21:03

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U F2 | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 74 | | 40-139 |

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 194508

SDG No.:

Instrument ID: AFID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 08/24/2015 13:59 Calibration End Date: 08/24/2015 17:07 Calibration ID: 30361

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-194508/7 | A0082107.D |
| Level 2 | STD2 240-194508/8 | A0082108.D |
| Level 3 | STD3 240-194508/9 | A0082109.D |
| Level 4 | STD4 240-194508/10 | A0082110.D |
| Level 5 | STD5 240-194508/11 | A0082111.D |
| Level 6 | STD6 240-194508/12 | A0082112.D |

| ANALYTE | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 6 | | | | | RT WINDOW | AVG RT |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 11.460 | 11.460 | 11.460 | 11.460 | 11.460 | 11.460 | | | | | 5.807 - 17.114 | 11.460 |
| C6-C12 | 12.725 | 12.725 | 12.725 | 12.725 | 12.725 | 12.725 | | | | | 5.807 - 19.644 | 12.725 |
| Trifluorotoluene (Surr) | 10.344 | 10.348 | 10.347 | 10.341 | 10.339 | 10.339 | | | | | 10.290 - 10.390 | 10.343 |

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 194508

SDG No.:

Instrument ID: AFID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 08/24/2015 13:59 Calibration End Date: 08/24/2015 17:07 Calibration ID: 30361

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-194508/7 | A0082107.D |
| Level 2 | STD2 240-194508/8 | A0082108.D |
| Level 3 | STD3 240-194508/9 | A0082109.D |
| Level 4 | STD4 240-194508/10 | A0082110.D |
| Level 5 | STD5 240-194508/11 | A0082111.D |
| Level 6 | STD6 240-194508/12 | A0082112.D |

| ANALYTE | CF | | | | CURVE TYPE | COEFFICIENT | | | # | MIN CF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 5 | LVL 2<br>LVL 6 | LVL 3 | LVL 4 | | B | M1 | M2 | | | | | | | | |
| Gasoline Range Organics [C6 - C10] | 281051<br>441376 | 404289<br>439409 | 445616 | 417800 | Ave | | 404923.648 | | | | 15.5 | | 20.0 | | | |
| C6-C12 | 344884<br>484706 | 442549<br>478304 | 494744 | 455788 | Ave | | 450162.318 | | | | 12.2 | | 20.0 | | | |
| Trifluorotoluene (Surr) | 452208<br>466474 | 509634<br>371094 | 436510 | 429984 | Ave | | 444317.252 | | | | 10.3 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 194508

SDG No.:

Instrument ID: AFID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 08/24/2015 13:59 Calibration End Date: 08/24/2015 17:07 Calibration ID: 30361

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-194508/7 | A0082107.D |
| Level 2 | STD2 240-194508/8 | A0082108.D |
| Level 3 | STD3 240-194508/9 | A0082109.D |
| Level 4 | STD4 240-194508/10 | A0082110.D |
| Level 5 | STD5 240-194508/11 | A0082111.D |
| Level 6 | STD6 240-194508/12 | A0082112.D |

| ANALYTE | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|
| | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Gasoline Range Organics [C6 - C10] | Ave | 28105105<br>703054501 | 80857808 | 178246588 | 334240056 | 529651436 | 100<br>1600 | 200 | 400 | 800 | 1200 |
| C6-C12 | Ave | 34488374<br>765287082 | 88509751 | 197897550 | 364630076 | 581646620 | 100<br>1600 | 200 | 400 | 800 | 1200 |
| Trifluorotoluene (Surr) | Ave | 452208<br>18554703 | 1019267 | 2182548 | 4299843 | 9329481 | 1.00<br>50.0 | 2.00 | 5.00 | 10.0 | 20.0 |

Curve Type Legend:
Ave = Average

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-194508/13 Calibration Date: 08/24/2015 17:45

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0082113.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 425225 | | 840 | 800 | 5.0 | 20.0 |
| C6-C12 | Ave | 450162 | 466294 | | 829 | 800 | 3.6 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 487053 | | 11.0 | 10.0 | 9.6 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-194508/13 Calibration Date: 08/24/2015 17:45

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0082113.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.46 | 5.81 | 17.11 |
| C6-C12 | 12.73 | 5.81 | 19.64 |
| Trifluorotoluene (Surr) | 10.34 | 10.29 | 10.39 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVRT 240-206211/4 Calibration Date: 11/11/2015 08:08

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111104.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 420445 | | 831 | 800 | 3.8 | 20.0 |
| C6-C12 | Ave | 450162 | 459766 | | 817 | 800 | 2.1 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 408705 | | 9.20 | 10.0 | -8.0 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVRT 240-206211/4 Calibration Date: 11/11/2015 08:08

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111104.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.54 | 5.94 | 17.18 |
| C6-C12 | 12.80 | 5.94 | 19.70 |
| Trifluorotoluene (Surr) | 10.36 | 10.31 | 10.41 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206211/6 Calibration Date: 11/11/2015 09:20

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111106.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 428576 | | 847 | 800 | 5.8 | 20.0 |
| C6-C12 | Ave | 450162 | 466289 | | 829 | 800 | 3.6 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 402849 | | 9.07 | 10.0 | -9.3 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206211/6 Calibration Date: 11/11/2015 09:20

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111106.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.54 | 5.94 | 17.18 |
| C6-C12 | 12.80 | 5.94 | 19.70 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206211/18 Calibration Date: 11/11/2015 16:42

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111118.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 421224 | | 832 | 800 | 4.0 | 20.0 |
| C6-C12 | Ave | 450162 | 459626 | | 817 | 800 | 2.1 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 402080 | | 9.05 | 10.0 | -9.5 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206211/18 Calibration Date: 11/11/2015 16:42

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111118.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.56 | 5.94 | 17.18 |
| C6-C12 | 12.82 | 5.94 | 19.70 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206211/28 Calibration Date: 11/11/2015 22:54

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111128.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 417296 | | 824 | 800 | 3.1 | 20.0 |
| C6-C12 | Ave | 450162 | 454853 | | 808 | 800 | 1.0 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 401488 | | 9.04 | 10.0 | -9.6 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206211/28 Calibration Date: 11/11/2015 22:54

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111128.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.56 | 5.94 | 17.18 |
| C6-C12 | 12.82 | 5.94 | 19.70 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVRT 240-206404/4 Calibration Date: 11/12/2015 07:26

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111204.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 427634 | | 845 | 800 | 5.6 | 20.0 |
| C6-C12 | Ave | 450162 | 465474 | | 827 | 800 | 3.4 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 405884 | | 9.14 | 10.0 | -8.6 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVRT 240-206404/4 Calibration Date: 11/12/2015 07:26

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111204.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.53 | 5.93 | 17.14 |
| C6-C12 | 12.79 | 5.93 | 19.66 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206404/8  Calibration Date: 11/12/2015 09:54

Instrument ID: AFID  Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX  ID: 0.45(mm)  Calib End Date: 08/24/2015 17:07

Lab File ID: A0111208.D  Conc. Units: ug/L  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 428972 | | 848 | 800 | 5.9 | 20.0 |
| C6-C12 | Ave | 450162 | 465297 | | 827 | 800 | 3.4 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 420936 | | 9.47 | 10.0 | -5.3 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206404/8 Calibration Date: 11/12/2015 09:54

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111208.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.53 | 5.93 | 17.14 |
| C6-C12 | 12.79 | 5.93 | 19.66 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-206227/1-A

Matrix: Solid  Lab File ID: A0111107.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/11/2015 09:56

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206211  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 46 | U | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 73 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-206444/1-A

Matrix: Solid  Lab File ID: A0111205.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/12/2015 08:03

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206404  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 46 | U | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 72 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-206227/2-A

Matrix: Solid  Lab File ID: A0111108.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/11/2015 10:31

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206211  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 786 | | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 95 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-206444/2-A

Matrix: Solid  Lab File ID: A0111206.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/12/2015 08:40

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206404  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 787 | | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 100 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: A0111126.D

Analysis Method: 8015C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.04(g) | Date Analyzed: 11/11/2015 21:40

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 856 | | 120 | 53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 99 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD | Lab Sample ID: 240-57597-15 MSD

Matrix: Solid | Lab File ID: A0111127.D

Analysis Method: 8015C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.05(g) | Date Analyzed: 11/11/2015 22:17

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 206 | | 120 | 53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 40 | | 40-139 |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: AFID Start Date: 08/24/2015 13:59

Analysis Batch Number: 194508 End Date: 08/24/2015 17:45

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-194508/7 IC | | 08/24/2015 13:59 | 1 | A0082107.D | RTX-VRX 0.45(mm) |
| STD2 240-194508/8 IC | | 08/24/2015 14:37 | 1 | A0082108.D | RTX-VRX 0.45(mm) |
| STD3 240-194508/9 IC | | 08/24/2015 15:14 | 1 | A0082109.D | RTX-VRX 0.45(mm) |
| STD4 240-194508/10 ICRT | | 08/24/2015 15:53 | 1 | A0082110.D | RTX-VRX 0.45(mm) |
| STD5 240-194508/11 IC | | 08/24/2015 16:31 | 1 | A0082111.D | RTX-VRX 0.45(mm) |
| STD6 240-194508/12 IC | | 08/24/2015 17:07 | 1 | A0082112.D | RTX-VRX 0.45(mm) |
| ICV 240-194508/13 | | 08/24/2015 17:45 | 1 | A0082113.D | RTX-VRX 0.45(mm) |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: AFID Start Date: 11/11/2015 08:08

Analysis Batch Number: 206211 End Date: 11/11/2015 22:54

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCVRT 240-206211/4 | | 11/11/2015 08:08 | 1 | A0111104.D | RTX-VRX 0.45(mm) |
| CCV 240-206211/6 | | 11/11/2015 09:20 | 1 | A0111106.D | RTX-VRX 0.45(mm) |
| MB 240-206227/1-A | | 11/11/2015 09:56 | 1 | A0111107.D | RTX-VRX 0.45(mm) |
| LCS 240-206227/2-A | | 11/11/2015 10:31 | 1 | A0111108.D | RTX-VRX 0.45(mm) |
| 240-57597-1 | F-VB-UST-1-14 | 11/11/2015 11:08 | 1 | A0111109.D | RTX-VRX 0.45(mm) |
| 240-57597-2 | F-VB-UST-2-14 | 11/11/2015 11:45 | 1 | A0111110.D | RTX-VRX 0.45(mm) |
| 240-57597-3 | F-VB-UST-3-14 | 11/11/2015 12:23 | 1 | A0111111.D | RTX-VRX 0.45(mm) |
| 240-57597-4 | F-VB-UST-4-14 | 11/11/2015 13:00 | 1 | A0111112.D | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/11/2015 13:36 | 1 | | RTX-VRX 0.45(mm) |
| 240-57597-6 | F-VB-UST-6-14 | 11/11/2015 14:13 | 1 | A0111114.D | RTX-VRX 0.45(mm) |
| 240-57597-7 | F-VB-UST-7-14 | 11/11/2015 14:50 | 1 | A0111115.D | RTX-VRX 0.45(mm) |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 15:27 | 1 | A0111116.D | RTX-VRX 0.45(mm) |
| CCV 240-206211/18 | | 11/11/2015 16:42 | 1 | A0111118.D | RTX-VRX 0.45(mm) |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 17:19 | 1 | A0111119.D | RTX-VRX 0.45(mm) |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 17:57 | 1 | A0111120.D | RTX-VRX 0.45(mm) |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 18:34 | 1 | A0111121.D | RTX-VRX 0.45(mm) |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 19:12 | 1 | A0111122.D | RTX-VRX 0.45(mm) |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 19:48 | 1 | A0111123.D | RTX-VRX 0.45(mm) |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 20:25 | 1 | A0111124.D | RTX-VRX 0.45(mm) |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 21:03 | 1 | A0111125.D | RTX-VRX 0.45(mm) |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 21:40 | 1 | A0111126.D | RTX-VRX 0.45(mm) |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 22:17 | 1 | A0111127.D | RTX-VRX 0.45(mm) |
| CCV 240-206211/28 | | 11/11/2015 22:54 | 1 | A0111128.D | RTX-VRX 0.45(mm) |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: AFID Start Date: 11/12/2015 07:26

Analysis Batch Number: 206404 End Date: 11/13/2015 01:22

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCVRT 240-206404/4 | | 11/12/2015 07:26 | 1 | A0111204.D | RTX-VRX 0.45(mm) |
| MB 240-206444/1-A | | 11/12/2015 08:03 | 1 | A0111205.D | RTX-VRX 0.45(mm) |
| LCS 240-206444/2-A | | 11/12/2015 08:40 | 1 | A0111206.D | RTX-VRX 0.45(mm) |
| 240-57597-5 | F-VB-UST-5-14 | 11/12/2015 09:17 | 1 | A0111207.D | RTX-VRX 0.45(mm) |
| CCV 240-206404/8 | | 11/12/2015 09:54 | 1 | A0111208.D | RTX-VRX 0.45(mm) |
| CCV 240-206404/16 | | 11/12/2015 14:54 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/12/2015 15:31 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/12/2015 16:09 | 1 | | RTX-VRX 0.45(mm) |
| CCV 240-206404/22 | | 11/12/2015 18:38 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/12/2015 19:51 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/12/2015 20:28 | 1 | | RTX-VRX 0.45(mm) |
| CCV 240-206404/33 | | 11/13/2015 01:22 | 1 | | RTX-VRX 0.45(mm) |

# GC VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 206227  Batch Start Date: 11/11/15 06:58  Batch Analyst: Reed, Rachel T

Batch Method: 5030A  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| MB 240-206227/1 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| LCS 240-206227/2 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| 240-57597-C-1 | F-VB-UST-1-14 | 5030A, 8015C | T | 4.98 g | 5 mL | | | | |
| 240-57597-C-2 | F-VB-UST-2-14 | 5030A, 8015C | T | 4.98 g | 5 mL | | | | |
| 240-57597-C-3 | F-VB-UST-3-14 | 5030A, 8015C | T | 4.99 g | 5 mL | | | | |
| 240-57597-C-4 | F-VB-UST-4-14 | 5030A, 8015C | T | 4.99 g | 5 mL | | | | |
| 240-57597-C-6 | F-VB-UST-6-14 | 5030A, 8015C | T | 5.01 g | 5 mL | | | | |
| 240-57597-C-7 | F-VB-UST-7-14 | 5030A, 8015C | T | 4.97 g | 5 mL | | | | |
| 240-57597-C-8 | F-VB-UST-8-14 | 5030A, 8015C | T | 4.98 g | 5 mL | | | | |
| 240-57597-C-9 | F-VB-UST-9-14 | 5030A, 8015C | T | 5.00 g | 5 mL | | | | |
| 240-57597-C-10 | F-VB-UST-10-14 | 5030A, 8015C | T | 5.02 g | 5 mL | | | | |
| 240-57597-C-11 | F-VB-UST-11-14 | 5030A, 8015C | T | 5.03 g | 5 mL | | | | |
| 240-57597-C-12 | F-VB-UST-12-14 | 5030A, 8015C | T | 5.04 g | 5 mL | | | | |
| 240-57597-C-13 | F-VB-UST-13-14 | 5030A, 8015C | T | 4.99 g | 5 mL | | | | |
| 240-57597-C-14 | F-VB-UST-14-14 | 5030A, 8015C | T | 5.02 g | 5 mL | | | | |
| 240-57597-C-15 | F-VB-UST-DUP | 5030A, 8015C | T | 5.00 g | 5 mL | | | | |
| 240-57597-C-15 MS | F-VB-UST-DUP | 5030A, 8015C | T | 5.04 g | 5 mL | | | | |
| 240-57597-C-15 MSD | F-VB-UST-DUP | 5030A, 8015C | T | 5.05 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| Balance ID | B045 |
| Blank Soil Lot Number | 1757979 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 206444  Batch Start Date: 11/12/15 07:59  Batch Analyst: Reed, Rachel T

Batch Method: 5030A  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| MB 240-206444/1 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| LCS 240-206444/2 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| 240-57597-C-5 | F-VB-UST-5-14 | 5030A, 8015C | T | 4.98 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| Balance ID | B045 |
| Blank Soil Lot Number | 1757979 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8015C - DRO

## Nonhalogenated Organics using GC/FID - Modified (Diesel Range Organics) by Method 8015C

FORM II
DIESEL RANGE ORGANICS SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): ID:

| Client Sample ID | Lab Sample ID | OTPH1 # |
|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 62 |
| F-VB-UST-2-14 | 240-57597-2 | 62 |
| F-VB-UST-3-14 | 240-57597-3 | 59 |
| F-VB-UST-4-14 | 240-57597-4 | 60 |
| F-VB-UST-5-14 | 240-57597-5 | 56 |
| F-VB-UST-6-14 | 240-57597-6 | 60 |
| F-VB-UST-7-14 | 240-57597-7 | 65 |
| F-VB-UST-8-14 | 240-57597-8 | 65 |
| F-VB-UST-9-14 | 240-57597-9 | 57 |
| F-VB-UST-10-14 | 240-57597-10 | 64 |
| F-VB-UST-11-14 | 240-57597-11 | 64 |
| F-VB-UST-12-14 | 240-57597-12 | 73 |
| F-VB-UST-13-14 | 240-57597-13 | 68 |
| F-VB-UST-14-14 | 240-57597-14 | 61 |
| F-VB-UST-DUP | 240-57597-15 | 65 |
| | MB 240-205788/22-A | 56 |
| | LCS 240-205788/23-A | 67 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 81 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 73 |

| | QC LIMITS |
|---|---|
| OTPH = o-Terphenyl (Surr) | 40-160 |

# Column to be used to flag recovery values

FORM II 8015C

FORM III
DIESEL RANGE ORGANICS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11009.D

Lab ID: LCS 240-205788/23-A Client ID:

| COMPOUND | SPIKE ADDED (mg/Kg) | LCS CONCENTRATION (mg/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Diesel | 84.0 | 53.7 | 64 | 52-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
DIESEL RANGE ORGANICS MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11113.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (mg/Kg) | SAMPLE CONCENTRATION (mg/Kg) | MS CONCENTRATION (mg/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Diesel | 99.5 | 7.5 U | 75.1 | 75 | 10-124 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
DIESEL RANGE ORGANICS MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11114.D

Lab ID: 240-57597-15 MSD Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (mg/Kg) | MSD CONCENTRATION (mg/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Diesel | 97.6 | 77.9 | 80 | 4 | 40 | 10-124 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM IV
DIESEL RANGE ORGANICS METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: MB 240-205788/22-A

Matrix: Solid  Date Extracted: 11/09/2015 09:40

Lab File ID:(1) P5B11008.D  Lab File ID:(2)

Date Analyzed:(1) 11/10/2015 16:59  Date Analyzed:(2)

Instrument ID:(1) A2HP5R  Instrument ID:(2)

GC Column:(1)  ID:  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-205788/23-A | 11/10/2015 17:31 | |
| F-VB-UST-1-14 | 240-57597-1 | 11/10/2015 20:07 | |
| F-VB-UST-2-14 | 240-57597-2 | 11/10/2015 20:38 | |
| F-VB-UST-3-14 | 240-57597-3 | 11/10/2015 21:10 | |
| F-VB-UST-4-14 | 240-57597-4 | 11/10/2015 21:41 | |
| F-VB-UST-5-14 | 240-57597-5 | 11/10/2015 22:43 | |
| F-VB-UST-6-14 | 240-57597-6 | 11/10/2015 23:14 | |
| F-VB-UST-7-14 | 240-57597-7 | 11/10/2015 23:46 | |
| F-VB-UST-8-14 | 240-57597-8 | 11/11/2015 00:17 | |
| F-VB-UST-9-14 | 240-57597-9 | 11/11/2015 00:48 | |
| F-VB-UST-10-14 | 240-57597-10 | 11/11/2015 01:19 | |
| F-VB-UST-11-14 | 240-57597-11 | 11/11/2015 01:50 | |
| F-VB-UST-12-14 | 240-57597-12 | 11/11/2015 19:03 | |
| F-VB-UST-13-14 | 240-57597-13 | 11/11/2015 19:34 | |
| F-VB-UST-14-14 | 240-57597-14 | 11/11/2015 20:06 | |
| F-VB-UST-DUP | 240-57597-15 | 11/11/2015 20:37 | |
| F-VB-UST-DUP MS | 240-57597-15 MS | 11/11/2015 21:08 | |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 11/11/2015 21:39 | |

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Sample No.: STD3 240-205608/10 Date Analyzed: 11/06/2015 17:59

Instrument ID: A2HP5R GC Column: ID: ()

Lab File ID (Standard): P5B10610.D Heated Purge: (Y/N) N

Calibration ID: 31386

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| INITIAL CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| STD3 240-205608/10 ICRT | | 11/06/2015 17:59 | P5B10610.D | 9.83 | | |

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206105/7 Date Analyzed: 11/10/2015 16:28

Instrument ID: A2HP5R GC Column: ID: ()

Lab File ID (Standard): P5B11007.D Heated Purge: (Y/N) N

Calibration ID: 31386

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCV 240-206105/7 CCVRT | | 11/10/2015 16:28 | P5B11007.D | 9.83 | | |
| MB 240-205788/22-A | | 11/10/2015 16:59 | P5B11008.D | 9.83 | | |
| LCS 240-205788/23-A | | 11/10/2015 17:31 | P5B11009.D | 9.83 | | |
| 240-57597-1 | F-VB-UST-1-14 | 11/10/2015 20:07 | P5B11014.D | 9.83 | | |
| 240-57597-2 | F-VB-UST-2-14 | 11/10/2015 20:38 | P5B11015.D | 9.83 | | |
| 240-57597-3 | F-VB-UST-3-14 | 11/10/2015 21:10 | P5B11016.D | 9.83 | | |
| 240-57597-4 | F-VB-UST-4-14 | 11/10/2015 21:41 | P5B11017.D | 9.83 | | |
| CCV 240-206105/18 | | 11/10/2015 22:12 | P5B11018.D | 9.83 | | |
| 240-57597-5 | F-VB-UST-5-14 | 11/10/2015 22:43 | P5B11019.D | 9.83 | | |
| 240-57597-6 | F-VB-UST-6-14 | 11/10/2015 23:14 | P5B11020.D | 9.83 | | |
| 240-57597-7 | F-VB-UST-7-14 | 11/10/2015 23:46 | P5B11021.D | 9.83 | | |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 00:17 | P5B11022.D | 9.83 | | |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 00:48 | P5B11023.D | 9.83 | | |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 01:19 | P5B11024.D | 9.83 | | |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 01:50 | P5B11025.D | 9.83 | | |
| CCV 240-206105/30 | | 11/11/2015 04:56 | P5B11031.D | 9.83 | | |

OTPH = o-Terphenyl

OTPH RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206356/7 Date Analyzed: 11/11/2015 18:00

Instrument ID: A2HP5R GC Column: ID: ()

Lab File ID (Standard): P5B11107.D Heated Purge: (Y/N) N

Calibration ID: 31386

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCV 240-206356/7 CCVRT | | 11/11/2015 18:00 | P5B11107.D | 9.83 | | |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 19:03 | P5B11109.D | 9.83 | | |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 19:34 | P5B11110.D | 9.83 | | |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 20:06 | P5B11111.D | 9.83 | | |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 20:37 | P5B11112.D | 9.83 | | |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 21:08 | P5B11113.D | 9.83 | | |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 21:39 | P5B11114.D | 9.83 | | |
| CCV 240-206356/15 | | 11/11/2015 22:11 | P5B11115.D | 9.83 | | |

OTPH = o-Terphenyl

OTPH RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: P5B11014.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.49(g) | Date Analyzed: 11/10/2015 20:07

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 16 | J B | 19 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 62 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: P5B11015.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:02

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.69(g) | Date Analyzed: 11/10/2015 20:38

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 10 | J B | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 62 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: P5B11016.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:04

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.32(g) | Date Analyzed: 11/10/2015 21:10

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 17.8 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.7 | U | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 59 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: P5B11017.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:06

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.31(g) | Date Analyzed: 11/10/2015 21:41

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.6 | J B | 20 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 60 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: P5B11019.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:08

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.01(g) | Date Analyzed: 11/10/2015 22:43

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.5 | U | 20 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 56 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: P5B11020.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:10

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.79(g) | Date Analyzed: 11/10/2015 23:14

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.7 | U | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 60 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: P5B11021.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:12

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.99(g) | Date Analyzed: 11/10/2015 23:46

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 16 | J B | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 65 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: P5B11022.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:14

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.45(g) | Date Analyzed: 11/11/2015 00:17

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 18.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.7 | U | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 65 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: P5B11023.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:16

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.38(g) | Date Analyzed: 11/11/2015 00:48

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 19.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.8 | U | 20 | 7.8 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 57 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: P5B11024.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:18

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.04(g) | Date Analyzed: 11/11/2015 01:19

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 13.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 13 | J B | 19 | 7.4 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 64 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: P5B11025.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:20

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.60(g) | Date Analyzed: 11/11/2015 01:50

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 9.0 | J B | 20 | 7.8 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 64 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: P5B11109.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:22

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.52(g) | Date Analyzed: 11/11/2015 19:03

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 30 | B | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 73 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: P5B11110.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:24

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.73(g) | Date Analyzed: 11/11/2015 19:34

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 14.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 13 | J B | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 68 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: P5B11111.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:26

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.43(g) | Date Analyzed: 11/11/2015 20:06

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 11 | J B | 19 | 7.4 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 61 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: P5B11112.D

Analysis Method: 8015C | Date Collected: 11/05/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.98(g) | Date Analyzed: 11/11/2015 20:37

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 8.2 | J B | 20 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 65 | | 40-160 |

FORM VI
DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205608

SDG No.:

Instrument ID: A2HP5R GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 11/06/2015 15:50 Calibration End Date: 11/06/2015 19:01 Calibration ID: 31386

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-205608/8 | P5B10608.D |
| Level 2 | STD2 240-205608/9 | P5B10609.D |
| Level 3 | STD3 240-205608/10 | P5B10610.D |
| Level 4 | STD4 240-205608/11 | P5B10611.D |
| Level 5 | STD5 240-205608/12 | P5B10612.D |

| ANALYTE | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | | | | | RT WINDOW | AVG RT |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | 9.331 | 9.331 | 9.331 | 9.331 | 9.331 | | | | | | 5.131 - 13.532 | 9.331 |
| o-Terphenyl (Surr) | 9.828 | 9.831 | 9.829 | 9.830 | 9.833 | | | | | | 9.728 - 9.928 | 9.830 |

FORM VI
DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205608

SDG No.:

Instrument ID: A2HP5R GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 11/06/2015 15:50 Calibration End Date: 11/06/2015 19:01 Calibration ID: 31386

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-205608/8 | P5B10608.D |
| Level 2 | STD2 240-205608/9 | P5B10609.D |
| Level 3 | STD3 240-205608/10 | P5B10610.D |
| Level 4 | STD4 240-205608/11 | P5B10611.D |
| Level 5 | STD5 240-205608/12 | P5B10612.D |

| ANALYTE | CF | | | | CURVE TYPE | COEFFICIENT | | | # | MIN CF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 5 | LVL 2 | LVL 3 | LVL 4 | | B | M1 | M2 | | | | | | | | |
| Diesel Range Organics [C10 - C28] | 156067 174183 | 161462 | 135888 | 139742 | Ave | | 153468.493 | | | | 10.3 | | 20.0 | | | |
| o-Terphenyl (Surr) | 213204 240558 | 209043 | 182141 | 189552 | Ave | | 206899.578 | | | | 11.1 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205608

SDG No.:

Instrument ID: A2HP5R GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 11/06/2015 15:50 Calibration End Date: 11/06/2015 19:01 Calibration ID: 31386

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-205608/8 | P5B10608.D |
| Level 2 | STD2 240-205608/9 | P5B10609.D |
| Level 3 | STD3 240-205608/10 | P5B10610.D |
| Level 4 | STD4 240-205608/11 | P5B10611.D |
| Level 5 | STD5 240-205608/12 | P5B10612.D |

| ANALYTE | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|
| | | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Diesel Range Organics [C10 - C28] | Ave | 15610573 | 32300497 | 67960720 | 139776482 | 348453260 | 100 | 200 | 500 | 1000 | 2000 |
| o-Terphenyl (Surr) | Ave | 1705635 | 3344683 | 7285634 | 15164183 | 38489230 | 8.00 | 16.0 | 40.0 | 80.0 | 160 |

Curve Type Legend:
Ave = Average

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-205608/13  Calibration Date: 11/06/2015 19:33

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B10613.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 164201 | | 537 | 502 | 7.0 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-205608/13 Calibration Date: 11/06/2015 19:33

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B10613.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206105/7  Calibration Date: 11/10/2015 16:28

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11007.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 149333 | | 487 | 500 | -2.7 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 178269 | | 34.5 | 40.0 | -13.8 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206105/7 Calibration Date: 11/10/2015 16:28

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11007.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206105/18  Calibration Date: 11/10/2015 22:12

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11018.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 166560 | | 543 | 500 | 8.5 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 216799 | | 41.9 | 40.0 | 4.8 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206105/18 Calibration Date: 11/10/2015 22:12

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11018.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206105/30  Calibration Date: 11/11/2015 04:56

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11031.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 176144 | | 574 | 500 | 14.8 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 226941 | | 43.9 | 40.0 | 9.7 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206105/30 Calibration Date: 11/11/2015 04:56

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11031.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206356/7  Calibration Date: 11/11/2015 18:00

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11107.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 156558 | | 510 | 500 | 2.0 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 215548 | | 41.7 | 40.0 | 4.2 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206356/7 Calibration Date: 11/11/2015 18:00

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11107.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206356/15  Calibration Date: 11/11/2015 22:11

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11115.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 166194 | | 542 | 500 | 8.3 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 218951 | | 42.3 | 40.0 | 5.8 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206356/15 Calibration Date: 11/11/2015 22:11

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11115.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205788/22-A

Matrix: Solid | Lab File ID: P5B11008.D

Analysis Method: 8015C | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30(g) | Date Analyzed: 11/10/2015 16:59

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.82 | J | 17 | 6.4 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 56 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-205788/23-A

Matrix: Solid | Lab File ID: P5B11009.D

Analysis Method: 8015C | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30(g) | Date Analyzed: 11/10/2015 17:31

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 67 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: P5B11113.D

Analysis Method: 8015C | Date Collected: 11/05/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.63(g) | Date Analyzed: 11/11/2015 21:08

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 77.0 | | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 81 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD | Lab Sample ID: 240-57597-15 MSD

Matrix: Solid | Lab File ID: P5B11114.D

Analysis Method: 8015C | Date Collected: 11/05/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.20(g) | Date Analyzed: 11/11/2015 21:39

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 79.5 | | 19 | 7.4 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 73 | | 40-160 |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP5R Start Date: 11/06/2015 15:50

Analysis Batch Number: 205608 End Date: 11/06/2015 19:33

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-205608/8 IC | | 11/06/2015 15:50 | 1 | P5B10608.D | |
| STD2 240-205608/9 IC | | 11/06/2015 17:28 | 1 | P5B10609.D | |
| STD3 240-205608/10 ICRT | | 11/06/2015 17:59 | 1 | P5B10610.D | |
| STD4 240-205608/11 IC | | 11/06/2015 18:30 | 1 | P5B10611.D | |
| STD5 240-205608/12 IC | | 11/06/2015 19:01 | 1 | P5B10612.D | |
| ICV 240-205608/13 | | 11/06/2015 19:33 | 1 | P5B10613.D | |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP5R Start Date: 11/10/2015 16:28

Analysis Batch Number: 206105 End Date: 11/11/2015 04:56

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-206105/7 CCVRT | | 11/10/2015 16:28 | 1 | P5B11007.D | |
| MB 240-205788/22-A | | 11/10/2015 16:59 | 1 | P5B11008.D | |
| LCS 240-205788/23-A | | 11/10/2015 17:31 | 1 | P5B11009.D | |
| ZZZZZ | | 11/10/2015 19:04 | 1 | | |
| 240-57597-1 | F-VB-UST-1-14 | 11/10/2015 20:07 | 1 | P5B11014.D | |
| 240-57597-2 | F-VB-UST-2-14 | 11/10/2015 20:38 | 1 | P5B11015.D | |
| 240-57597-3 | F-VB-UST-3-14 | 11/10/2015 21:10 | 1 | P5B11016.D | |
| 240-57597-4 | F-VB-UST-4-14 | 11/10/2015 21:41 | 1 | P5B11017.D | |
| CCV 240-206105/18 | | 11/10/2015 22:12 | 1 | P5B11018.D | |
| 240-57597-5 | F-VB-UST-5-14 | 11/10/2015 22:43 | 1 | P5B11019.D | |
| 240-57597-6 | F-VB-UST-6-14 | 11/10/2015 23:14 | 1 | P5B11020.D | |
| 240-57597-7 | F-VB-UST-7-14 | 11/10/2015 23:46 | 1 | P5B11021.D | |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 00:17 | 1 | P5B11022.D | |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 00:48 | 1 | P5B11023.D | |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 01:19 | 1 | P5B11024.D | |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 01:50 | 1 | P5B11025.D | |
| CCV 240-206105/30 | | 11/11/2015 04:56 | 1 | P5B11031.D | |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP5R Start Date: 11/11/2015 18:00

Analysis Batch Number: 206356 End Date: 11/11/2015 22:11

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-206356/7 CCVRT | | 11/11/2015 18:00 | 1 | P5B11107.D | |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 19:03 | 1 | P5B11109.D | |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 19:34 | 1 | P5B11110.D | |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 20:06 | 1 | P5B11111.D | |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 20:37 | 1 | P5B11112.D | |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 21:08 | 1 | P5B11113.D | |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 21:39 | 1 | P5B11114.D | |
| CCV 240-206356/15 | | 11/11/2015 22:11 | 1 | P5B11115.D | |

# DIESEL RANGE ORGANICS BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205788  Batch Start Date: 11/09/15 10:35  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/10/15 02:35

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exDIESELspkw 00068 | exOTPSURRW 00006 | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-D-1 | F-VB-UST-1-14 | 3540C, 8015C | T | 30.49 g | 5 mL | | 1 mL | | |
| 240-57597-E-2 | F-VB-UST-2-14 | 3540C, 8015C | T | 29.69 g | 5 mL | | 1 mL | | |
| 240-57597-D-3 | F-VB-UST-3-14 | 3540C, 8015C | T | 30.32 g | 5 mL | | 1 mL | | |
| 240-57597-D-4 | F-VB-UST-4-14 | 3540C, 8015C | T | 30.31 g | 5 mL | | 1 mL | | |
| 240-57597-D-5 | F-VB-UST-5-14 | 3540C, 8015C | T | 30.01 g | 5 mL | | 1 mL | | |
| 240-57597-D-6 | F-VB-UST-6-14 | 3540C, 8015C | T | 29.79 g | 5 mL | | 1 mL | | |
| 240-57597-D-7 | F-VB-UST-7-14 | 3540C, 8015C | T | 29.99 g | 5 mL | | 1 mL | | |
| 240-57597-E-8 | F-VB-UST-8-14 | 3540C, 8015C | T | 30.45 g | 5 mL | | 1 mL | | |
| 240-57597-D-9 | F-VB-UST-9-14 | 3540C, 8015C | T | 30.38 g | 5 mL | | 1 mL | | |
| 240-57597-D-10 | F-VB-UST-10-14 | 3540C, 8015C | T | 30.04 g | 5 mL | | 1 mL | | |
| 240-57597-E-11 | F-VB-UST-11-14 | 3540C, 8015C | T | 29.60 g | 5 mL | | 1 mL | | |
| 240-57597-E-12 | F-VB-UST-12-14 | 3540C, 8015C | T | 29.52 g | 5 mL | | 1 mL | | |
| 240-57597-E-13 | F-VB-UST-13-14 | 3540C, 8015C | T | 29.73 g | 5 mL | | 1 mL | | |
| 240-57597-D-14 | F-VB-UST-14-14 | 3540C, 8015C | T | 30.43 g | 5 mL | | 1 mL | | |
| 240-57597-D-15 | F-VB-UST-DUP | 3540C, 8015C | T | 29.98 g | 5 mL | | 1 mL | | |
| 240-57597-D-15 MS | F-VB-UST-DUP | 3540C, 8015C | T | 29.63 g | 5 mL | 1 mL | 1 mL | | |
| 240-57597-D-15 MSD | F-VB-UST-DUP | 3540C, 8015C | T | 30.20 g | 5 mL | 1 mL | 1 mL | | |
| MB 240-205788/22 | | 3540C, 8015C | | 30 g | 5 mL | | 1 mL | | |
| LCS 240-205788/23 | | 3540C, 8015C | | 30 g | 5 mL | 1 mL | 1 mL | | |

| Batch Notes | |
|---|---|
| Balance ID | 1339907 |
| Na2SO4 Lot Number | 2204509 |
| Prep Solvent Lot # | 2258960 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8082A

## Polychlorinated Biphenyls (PCBs) (GC) by Method 8082A

FORM II
PCBS SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): CLP-2 ID: 0.53(mm)

| Client Sample ID | Lab Sample ID | TCX1 # | DCB1 # |
|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 78 | 87 |
| F-VB-UST-2-14 | 240-57597-2 | 84 | 100 |
| F-VB-UST-3-14 | 240-57597-3 | 68 | 95 |
| F-VB-UST-4-14 | 240-57597-4 | 75 | 104 |
| F-VB-UST-5-14 | 240-57597-5 | 73 | 91 |
| F-VB-UST-6-14 | 240-57597-6 | 76 | 109 |
| F-VB-UST-7-14 | 240-57597-7 | 76 | 98 |
| F-VB-UST-8-14 | 240-57597-8 | 72 | 90 |
| F-VB-UST-9-14 | 240-57597-9 | 77 | 95 |
| F-VB-UST-10-14 | 240-57597-10 | 83 | 89 |
| F-VB-UST-11-14 | 240-57597-11 | 82 | 95 |
| F-VB-UST-12-14 | 240-57597-12 | 81 | 92 |
| F-VB-UST-13-14 | 240-57597-13 | 77 | 94 |
| F-VB-UST-14-14 | 240-57597-14 | 78 | 89 |
| F-VB-UST-DUP | 240-57597-15 | 74 | 89 |
| | MB 240-205809/21-A | 75 | 89 |
| | LCS 240-205809/22-A | 75 | 85 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 81 | 95 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 77 | 89 |

| | QC LIMITS |
|---|---|
| TCX = Tetrachloro-m-xylene | 13-134 |
| DCB = DCB Decachlorobiphenyl | 10-155 |

# Column to be used to flag recovery values

FORM II 8082A

FORM III
PCBS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P4000027.D

Lab ID: LCS 240-205809/22-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Aroclor-1016 | 333 | 196 | 59 | 51-120 | |
| Aroclor-1260 | 334 | 201 | 60 | 48-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM III
PCBS MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P4000030.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Aroclor-1016 | 394 | 14 U | 244 | 62 | 17-141 | |
| Aroclor-1260 | 395 | 10 U | 260 | 66 | 19-137 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM III
PCBS MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P4000031.D

Lab ID: 240-57597-15 MSD Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Aroclor-1016 | 393 | 229 | 58 | 9 | 56 | 17-141 | |
| Aroclor-1260 | 393 | 238 | 61 | 9 | 55 | 19-137 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM IV
PCBS METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: MB 240-205809/21-A

Matrix: Solid  Date Extracted: 11/09/2015 09:53

Lab File ID:(1) P4000026.D  Lab File ID:(2)

Date Analyzed:(1) 11/11/2015 17:17  Date Analyzed:(2)

Instrument ID:(1) A2HP4  Instrument ID:(2)

GC Column:(1) CLP-2  ID: 0.53(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 11/11/2015 13:28 | |
| F-VB-UST-2-14 | 240-57597-2 | 11/11/2015 13:45 | |
| F-VB-UST-3-14 | 240-57597-3 | 11/11/2015 14:01 | |
| F-VB-UST-4-14 | 240-57597-4 | 11/11/2015 14:17 | |
| F-VB-UST-5-14 | 240-57597-5 | 11/11/2015 14:33 | |
| F-VB-UST-6-14 | 240-57597-6 | 11/11/2015 14:50 | |
| F-VB-UST-7-14 | 240-57597-7 | 11/11/2015 15:06 | |
| F-VB-UST-8-14 | 240-57597-8 | 11/11/2015 15:23 | |
| F-VB-UST-9-14 | 240-57597-9 | 11/11/2015 15:39 | |
| F-VB-UST-10-14 | 240-57597-10 | 11/11/2015 15:55 | |
| F-VB-UST-11-14 | 240-57597-11 | 11/11/2015 16:12 | |
| F-VB-UST-12-14 | 240-57597-12 | 11/11/2015 16:28 | |
| F-VB-UST-13-14 | 240-57597-13 | 11/11/2015 16:44 | |
| F-VB-UST-14-14 | 240-57597-14 | 11/11/2015 17:01 | |
| | LCS 240-205809/22-A | 11/11/2015 17:33 | |
| F-VB-UST-DUP | 240-57597-15 | 11/11/2015 18:06 | |
| F-VB-UST-DUP MS | 240-57597-15 MS | 11/11/2015 18:22 | |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 11/11/2015 18:39 | |

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: STD05 240-205222/31  Date Analyzed: 11/04/2015  22:13

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)

Lab File ID (Standard): P4000031.D  Heated Purge: (Y/N) N

Calibration ID: 31286

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 138265983 | 2.08 | | | | |
| UPPER LIMIT | | 276531966 | 2.58 | | | | |
| LOWER LIMIT | | 69132992 | 1.58 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-205222/39 | | 151134439 | 2.09 | | | | |
| CCV 240-206283/3 CCVIS | | 120153607 | 2.07 | | | | |

BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206283/3  Date Analyzed: 11/11/2015 11:01

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)

Lab File ID (Standard): P4000003.D  Heated Purge: (Y/N) N

Calibration ID: 31286

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 120153607 | 2.07 | | | | |
| UPPER LIMIT | | 240307214 | 2.57 | | | | |
| LOWER LIMIT | | 60076804 | 1.57 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-206283/4 | | 117662635 | 2.07 | | | | |
| CCV 240-206283/5 | | 128931083 | 2.08 | | | | |
| CCV 240-206283/6 | | 129660063 | 2.07 | | | | |
| CCV 240-206283/7 | | 124415610 | 2.07 | | | | |
| 240-57597-1 | F-VB-UST-1-14 | 140229279 | 2.08 | | | | |
| 240-57597-2 | F-VB-UST-2-14 | 131329155 | 2.07 | | | | |
| 240-57597-3 | F-VB-UST-3-14 | 142270888 | 2.07 | | | | |
| 240-57597-4 | F-VB-UST-4-14 | 139052581 | 2.07 | | | | |
| 240-57597-5 | F-VB-UST-5-14 | 150647200 | 2.07 | | | | |
| 240-57597-6 | F-VB-UST-6-14 | 136592279 | 2.07 | | | | |
| 240-57597-7 | F-VB-UST-7-14 | 142766511 | 2.07 | | | | |
| 240-57597-8 | F-VB-UST-8-14 | 143364145 | 2.07 | | | | |
| 240-57597-9 | F-VB-UST-9-14 | 135564000 | 2.07 | | | | |
| 240-57597-10 | F-VB-UST-10-14 | 134509895 | 2.07 | | | | |
| 240-57597-11 | F-VB-UST-11-14 | 146836344 | 2.07 | | | | |
| 240-57597-12 | F-VB-UST-12-14 | 141041341 | 2.07 | | | | |
| 240-57597-13 | F-VB-UST-13-14 | 143471346 | 2.08 | | | | |
| 240-57597-14 | F-VB-UST-14-14 | 143338015 | 2.07 | | | | |
| MB 240-205809/21-A | | 150167483 | 2.07 | | | | |
| LCS 240-205809/22-A | | 138157065 | 2.07 | | | | |

BNB = 1-Bromo-2-nitrobenzene
BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVIS 240-206283/28  Date Analyzed: 11/11/2015 17:50

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)

Lab File ID (Standard): P4000028.D  Heated Purge: (Y/N) N

Calibration ID: 31286

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 124491700 | 2.08 | | | | |
| UPPER LIMIT | | 248983400 | 2.58 | | | | |
| LOWER LIMIT | | 62245850 | 1.58 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| 240-57597-15 | F-VB-UST-DUP | 155311076 | 2.08 | | | | |
| 240-57597-15 MS | F-VB-UST-DUP MS | 144577114 | 2.07 | | | | |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 162028723 | 2.07 | | | | |

BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM X
IDENTIFICATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS Lab Sample ID: 240-57597-15 MS

Instrument ID (1): A2HP4 Instrument ID (2):

Date Analyzed (1): 11/11/2015 18:22 Date Analyzed (2):

GC Column (1): CLP-2 ID: 0.53(mm) GC Column (2): ID:

| ANALYTE | COL | PEAK | RT | RT WINDOW | | CONCENTRATION | | |
|---|---|---|---|---|---|---|---|---|
| | | | | FROM | TO | PEAK | MEAN | |
| Aroclor-1016 | 2 | 1 | 4.50 | 4.48 | 4.54 | 244 | 244 | |
| | | 2 | 5.18 | 5.16 | 5.22 | 246 | | |
| | | 3 | 5.89 | 5.87 | 5.93 | 245 | | |
| | | 4 | 6.09 | 6.07 | 6.13 | 229 | | |
| | | 5 | 6.73 | 6.71 | 6.77 | 256 | | |
| Aroclor-1260 | 2 | 1 | 8.10 | 8.08 | 8.14 | 267 | 260 | |
| | | 2 | 8.30 | 8.28 | 8.34 | 249 | | |
| | | 3 | 8.62 | 8.59 | 8.65 | 240 | | |
| | | 4 | 9.16 | 9.14 | 9.20 | 270 | | |
| | | 5 | 9.47 | 9.44 | 9.50 | 272 | | |

FORM X
IDENTIFICATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD Lab Sample ID: 240-57597-15 MSD

Instrument ID (1): A2HP4 Instrument ID (2):

Date Analyzed (1): 11/11/2015 18:39 Date Analyzed (2):

GC Column (1): CLP-2 ID: 0.53(mm) GC Column (2): ID:

| ANALYTE | COL | PEAK | RT | RT WINDOW | | CONCENTRATION | | |
|---|---|---|---|---|---|---|---|---|
| | | | | FROM | TO | PEAK | MEAN | |
| Aroclor-1016 | 2 | 1 | 4.50 | 4.48 | 4.54 | 231 | 229 | |
| | | 2 | 5.18 | 5.16 | 5.22 | 232 | | |
| | | 3 | 5.89 | 5.87 | 5.93 | 232 | | |
| | | 4 | 6.10 | 6.07 | 6.13 | 215 | | |
| | | 5 | 6.74 | 6.71 | 6.77 | 234 | | |
| Aroclor-1260 | 2 | 1 | 8.10 | 8.08 | 8.14 | 248 | 238 | |
| | | 2 | 8.30 | 8.28 | 8.34 | 228 | | |
| | | 3 | 8.62 | 8.59 | 8.65 | 218 | | |
| | | 4 | 9.17 | 9.14 | 9.20 | 246 | | |
| | | 5 | 9.47 | 9.44 | 9.50 | 251 | | |

FORM X
IDENTIFICATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Client Sample ID: Lab Sample ID: LCS 240-205809/22-A

Instrument ID (1): A2HP4 Instrument ID (2):

Date Analyzed (1): 11/11/2015 17:33 Date Analyzed (2):

GC Column (1): CLP-2 ID: 0.53(mm) GC Column (2): ID:

| ANALYTE | COL | PEAK | RT | RT WINDOW | | CONCENTRATION | | |
|---|---|---|---|---|---|---|---|---|
| | | | | FROM | TO | PEAK | MEAN | |
| Aroclor-1016 | 2 | 1 | 4.49 | 4.46 | 4.52 | 195 | 196 | |
| | | 2 | 5.18 | 5.15 | 5.21 | 198 | | |
| | | 3 | 5.89 | 5.85 | 5.91 | 196 | | |
| | | 4 | 6.09 | 6.06 | 6.12 | 189 | | |
| | | 5 | 6.73 | 6.70 | 6.76 | 203 | | |
| Aroclor-1260 | 2 | 1 | 8.10 | 8.06 | 8.12 | 211 | 201 | |
| | | 2 | 8.30 | 8.27 | 8.33 | 195 | | |
| | | 3 | 8.62 | 8.58 | 8.64 | 184 | | |
| | | 4 | 9.17 | 9.13 | 9.19 | 207 | | |
| | | 5 | 9.47 | 9.43 | 9.49 | 208 | | |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: P4000012.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.54(g) | Date Analyzed: 11/11/2015 13:28

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 40 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.6 | U | 40 | 9.6 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 78 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 87 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: P4000013.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:02

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.07(g) | Date Analyzed: 11/11/2015 13:45

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 39 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.4 | U | 39 | 9.4 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 39 | 16 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 84 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 100 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: P4000014.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:04

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.14(g) | Date Analyzed: 11/11/2015 14:01

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 17.8 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 15 | U | 40 | 15 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.7 | U | 40 | 9.7 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 68 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 95 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: P4000015.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:06

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.54(g) | Date Analyzed: 11/11/2015 14:17

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 40 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.6 | U | 40 | 9.6 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 75 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 104 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: P4000016.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:08

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.48(g) | Date Analyzed: 11/11/2015 14:33

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 73 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 91 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: P4000017.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:10

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.87(g) | Date Analyzed: 11/11/2015 14:50

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 40 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.6 | U | 40 | 9.6 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 76 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 109 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: P4000018.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:12

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.04(g) | Date Analyzed: 11/11/2015 15:06

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 16.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 76 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 98 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: P4000019.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:14

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.74(g) | Date Analyzed: 11/11/2015 15:23

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 18.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 15 | U | 41 | 15 |
| 11104-28-2 | Aroclor-1221 | 20 | U | 41 | 20 |
| 11141-16-5 | Aroclor-1232 | 25 | U | 41 | 25 |
| 53469-21-9 | Aroclor-1242 | 14 | U | 41 | 14 |
| 12672-29-6 | Aroclor-1248 | 9.8 | U | 41 | 9.8 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 41 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 41 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 72 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 90 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: P4000020.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:16

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.07(g) | Date Analyzed: 11/11/2015 15:39

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 19.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 15 | U | 41 | 15 |
| 11104-28-2 | Aroclor-1221 | 20 | U | 41 | 20 |
| 11141-16-5 | Aroclor-1232 | 25 | U | 41 | 25 |
| 53469-21-9 | Aroclor-1242 | 14 | U | 41 | 14 |
| 12672-29-6 | Aroclor-1248 | 9.9 | U | 41 | 9.9 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 41 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 41 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 77 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 95 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: P4000021.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:18

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.97(g) | Date Analyzed: 11/11/2015 15:55

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 13.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 83 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 89 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: P4000022.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:20

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.28(g) | Date Analyzed: 11/11/2015 16:12

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 82 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 95 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: P4000023.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:22

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.29(g) | Date Analyzed: 11/11/2015 16:28

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.4 | U | 39 | 9.4 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 39 | 16 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 81 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 92 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: P4000024.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:24

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.41(g) | Date Analyzed: 11/11/2015 16:44

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 14.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 77 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 94 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: P4000025.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:26

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.19(g) | Date Analyzed: 11/11/2015 17:01

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 78 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 89 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: P4000029.D

Analysis Method: 8082A | Date Collected: 11/05/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.34(g) | Date Analyzed: 11/11/2015 18:06

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 18 | U | 38 | 18 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.2 | U | 38 | 9.2 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 74 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 89 | | 10-155 |

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4 GC Column: CLP-2 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 14:51 Calibration End Date: 11/04/2015 16:13 Calibration ID: 31262

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/4 | P4000004.D |
| Level 2 | STD01 240-205222/5 | P4000005.D |
| Level 3 | STD02 240-205222/6 | P4000006.D |
| Level 4 | STD05 240-205222/7 | P4000007.D |
| Level 5 | STD1 240-205222/8 | P4000008.D |
| Level 6 | STD2 240-205222/9 | P4000009.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1232 Peak 1 | 0.0146<br>0.0122 | 0.0144 | 0.0134 | 0.0134 | 0.0122 | Lin1 | 0.0002 | 0.0122 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1232 Peak 2 | 0.0119<br>0.0097 | 0.0118 | 0.0110 | 0.0109 | 0.0098 | Lin1 | 0.0002 | 0.0098 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 3 | 0.0217<br>0.0205 | 0.0220 | 0.0209 | 0.0216 | 0.0202 | Lin1 | 0.0001 | 0.0205 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1232 Peak 4 | 0.0095<br>0.0081 | 0.0100 | 0.0094 | 0.0089 | 0.0081 | Lin1 | 0.0001 | 0.0081 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 5 | 0.0072<br>0.0065 | 0.0075 | 0.0069 | 0.0070 | 0.0066 | Lin1 | 0.0001 | 0.0066 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1262 Peak 1 | 0.0260<br>0.0242 | 0.0274 | 0.0250 | 0.0248 | 0.0248 | Lin1 | 0.0002 | 0.0244 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 2 | 0.0377<br>0.0377 | 0.0413 | 0.0382 | 0.0381 | 0.0383 | Lin1 | 0.0001 | 0.0379 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 3 | 0.0311<br>0.0318 | 0.0344 | 0.0320 | 0.0318 | 0.0323 | Lin1 | 0 | 0.0319 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 4 | 0.0698<br>0.0769 | 0.0788 | 0.0741 | 0.0751 | 0.0769 | Lin1 | 0 | 0.0769 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 5 | 0.0515<br>0.0538 | 0.0573 | 0.0535 | 0.0535 | 0.0541 | Lin1 | 0 | 0.0539 | | | | | | | 1.0000 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4 GC Column: CLP-2 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 14:51 Calibration End Date: 11/04/2015 16:13 Calibration ID: 31262

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/4 | P4000004.D |
| Level 2 | STD01 240-205222/5 | P4000005.D |
| Level 3 | STD02 240-205222/6 | P4000006.D |
| Level 4 | STD05 240-205222/7 | P4000007.D |
| Level 5 | STD1 240-205222/8 | P4000008.D |
| Level 6 | STD2 240-205222/9 | P4000009.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1232 Peak 1 | BNB | Lin1 | 1933285<br>63941977 | 4095893 | 7303079 | 16914408 | 33770823 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 2 | BNB | Lin1 | 1578310<br>50725715 | 3372584 | 6002899 | 13706058 | 27055994 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 3 | BNB | Lin1 | 2876493<br>107658914 | 6259111 | 11374836 | 27178028 | 55762073 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 4 | BNB | Lin1 | 1257572<br>42453230 | 2846698 | 5080531 | 11255540 | 22469169 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 5 | BNB | Lin1 | 958391<br>34172989 | 2139622 | 3763738 | 8806970 | 18208852 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 1 | BNB | Lin1 | 3439151<br>127156329 | 7810498 | 13593351 | 31298082 | 68292376 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 2 | BNB | Lin1 | 4981768<br>197819034 | 11766832 | 20756002 | 47996867 | 105623105 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 3 | BNB | Lin1 | 4115030<br>166759669 | 9809163 | 17399332 | 40126368 | 89102815 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 4 | BNB | Lin1 | 9238963<br>403469407 | 22450345 | 40269720 | 94574833 | 212075612 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 5 | BNB | Lin1 | 6817657<br>282427060 | 16339189 | 29055665 | 67469104 | 149093104 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 16:30  Calibration End Date: 11/04/2015 17:51  Calibration ID: 31268

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/10 | P4000010.D |
| Level 2 | STD01 240-205222/11 | P4000011.D |
| Level 3 | STD02 240-205222/12 | P4000012.D |
| Level 4 | STD05 240-205222/13 | P4000013.D |
| Level 5 | STD1 240-205222/14 | P4000014.D |
| Level 6 | STD2 240-205222/15 | P4000015.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1242 Peak 1 | 0.0108 0.0082 | 0.0098 | 0.0093 | 0.0090 | 0.0085 | Lin1 | 0.0002 | 0.0083 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 2 | 0.0208 0.0159 | 0.0186 | 0.0181 | 0.0174 | 0.0164 | Lin1 | 0.0003 | 0.0160 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 3 | 0.0407 0.0367 | 0.0373 | 0.0369 | 0.0376 | 0.0368 | Lin1 | 0.0002 | 0.0367 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1242 Peak 4 | 0.0183 0.0142 | 0.0165 | 0.0153 | 0.0150 | 0.0144 | Lin1 | 0.0002 | 0.0142 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1242 Peak 5 | 0.0169 0.0109 | 0.0131 | 0.0124 | 0.0117 | 0.0111 | Lin1 | 0.0003 | 0.0108 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 1 | 0.1004 0.0910 | 0.0938 | 0.0918 | 0.0903 | 0.0899 | Lin1 | 0.0004 | 0.0902 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 2 | 0.0969 0.0876 | 0.0916 | 0.0907 | 0.0881 | 0.0870 | Lin1 | 0.0005 | 0.0871 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 3 | 0.0848 0.0764 | 0.0793 | 0.0781 | 0.0771 | 0.0756 | Lin1 | 0.0004 | 0.0759 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 4 | 0.0354 0.0308 | 0.0332 | 0.0323 | 0.0312 | 0.0308 | Lin1 | 0.0003 | 0.0306 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 5 | 0.2560 0.2513 | 0.2411 | 0.2398 | 0.2407 | 0.2403 | Lin1 | 0 | 0.2467 | | | | | | | 0.9990 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 16:30  Calibration End Date: 11/04/2015 17:51  Calibration ID: 31268

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/10 | P4000010.D |
| Level 2 | STD01 240-205222/11 | P4000011.D |
| Level 3 | STD02 240-205222/12 | P4000012.D |
| Level 4 | STD05 240-205222/13 | P4000013.D |
| Level 5 | STD1 240-205222/14 | P4000014.D |
| Level 6 | STD2 240-205222/15 | P4000015.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1242 Peak 1 | BNB | Lin1 | 1416279<br>42642341 | 2631282 | 4898596 | 12129968 | 22349639 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 2 | BNB | Lin1 | 2727967<br>82584170 | 5016168 | 9472105 | 23377100 | 43234168 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 3 | BNB | Lin1 | 5339536<br>190275331 | 10059287 | 19366944 | 50515067 | 97245652 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 4 | BNB | Lin1 | 2396365<br>73778193 | 4446179 | 8021924 | 20164058 | 37966389 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 5 | BNB | Lin1 | 2219925<br>56371385 | 3525488 | 6485701 | 15661664 | 29226139 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 1 | BNB | Lin1 | 13163622<br>472145731 | 25289564 | 48180973 | 121179071 | 237699554 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 2 | BNB | Lin1 | 12703541<br>454549189 | 24701408 | 47581502 | 118240573 | 229940017 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 3 | BNB | Lin1 | 11123762<br>396177954 | 21390831 | 40984813 | 103417772 | 199923731 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 4 | BNB | Lin1 | 4645327<br>159800157 | 8942838 | 16959213 | 41849329 | 81325789 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 5 | BNB | Lin1 | 33570153<br>1303248607 | 65017075 | 125845634 | 322956363 | 635404036 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4 GC Column: CLP-2 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 18:08 Calibration End Date: 11/04/2015 19:30 Calibration ID: 31274

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/16 | P4000016.D |
| Level 2 | STD01 240-205222/17 | P4000017.D |
| Level 3 | STD02 240-205222/18 | P4000018.D |
| Level 4 | STD05 240-205222/19 | P4000019.D |
| Level 5 | STD1 240-205222/20 | P4000020.D |
| Level 6 | STD2 240-205222/21 | P4000021.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1248 Peak 1 | 0.0262 0.0244 | 0.0257 | 0.0252 | 0.0241 | 0.0238 | Lin1 | 0.0001 | 0.0241 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 2 | 0.0147 0.0119 | 0.0141 | 0.0133 | 0.0123 | 0.0118 | Lin1 | 0.0002 | 0.0118 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 3 | 0.0221 0.0182 | 0.0209 | 0.0200 | 0.0185 | 0.0178 | Lin1 | 0.0003 | 0.0180 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 4 | 0.0230 0.0230 | 0.0209 | 0.0242 | 0.0230 | 0.0224 | Lin1 | 0 | 0.0229 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1248 Peak 5 | 0.0111 0.0116 | 0.0121 | 0.0125 | 0.0118 | 0.0111 | Lin1 | 0 | 0.0115 | | | | | | | 0.9990 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4 GC Column: CLP-2 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 18:08 Calibration End Date: 11/04/2015 19:30 Calibration ID: 31274

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/16 | P4000016.D |
| Level 2 | STD01 240-205222/17 | P4000017.D |
| Level 3 | STD02 240-205222/18 | P4000018.D |
| Level 4 | STD05 240-205222/19 | P4000019.D |
| Level 5 | STD1 240-205222/20 | P4000020.D |
| Level 6 | STD2 240-205222/21 | P4000021.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1248 Peak 1 | BNB | Lin1 | 3621408<br>135686827 | 7056408 | 13893180 | 32542109 | 61737564 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 2 | BNB | Lin1 | 2026547<br>66471349 | 3863774 | 7340597 | 16627224 | 30666654 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 3 | BNB | Lin1 | 3052634<br>101195767 | 5751628 | 11041531 | 25008923 | 46323988 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 4 | BNB | Lin1 | 3174691<br>128291777 | 5746638 | 13346231 | 31077384 | 58123141 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 5 | BNB | Lin1 | 1538217<br>64509725 | 3319666 | 6884267 | 15873534 | 28832671 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 19:46  Calibration End Date: 11/04/2015 21:08  Calibration ID: 31280

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/22 | P4000022.D |
| Level 2 | STD01 240-205222/23 | P4000023.D |
| Level 3 | STD02 240-205222/24 | P4000024.D |
| Level 4 | STD05 240-205222/25 | P4000025.D |
| Level 5 | STD1 240-205222/26 | P4000026.D |
| Level 6 | STD2 240-205222/27 | P4000027.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1221 Peak 1 | 0.0082 0.0059 | 0.0071 | 0.0067 | 0.0065 | 0.0060 | Lin1 | 0.0001 | 0.0059 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1221 Peak 2 | 0.0056 0.0038 | 0.0048 | 0.0046 | 0.0044 | 0.0040 | Lin1 | 0.0001 | 0.0039 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1221 Peak 3 | 0.0205 0.0147 | 0.0173 | 0.0165 | 0.0161 | 0.0151 | Lin1 | 0.0003 | 0.0148 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1254 Peak 1 | 0.0288 0.0211 | 0.0242 | 0.0224 | 0.0222 | 0.0212 | Lin1 | 0.0004 | 0.0209 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 2 | 0.0430 0.0354 | 0.0364 | 0.0351 | 0.0362 | 0.0354 | Lin1 | 0.0003 | 0.0352 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 3 | 0.0328 0.0255 | 0.0276 | 0.0256 | 0.0267 | 0.0261 | Lin1 | 0.0003 | 0.0255 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 4 | 0.0251 0.0215 | 0.0209 | 0.0202 | 0.0214 | 0.0214 | Lin1 | 0.0001 | 0.0213 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1254 Peak 5 | 0.0446 0.0346 | 0.0372 | 0.0350 | 0.0360 | 0.0351 | Lin1 | 0.0004 | 0.0345 | | | | | | | 1.0000 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4 GC Column: CLP-2 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 19:46 Calibration End Date: 11/04/2015 21:08 Calibration ID: 31280

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/22 | P4000022.D |
| Level 2 | STD01 240-205222/23 | P4000023.D |
| Level 3 | STD02 240-205222/24 | P4000024.D |
| Level 4 | STD05 240-205222/25 | P4000025.D |
| Level 5 | STD1 240-205222/26 | P4000026.D |
| Level 6 | STD2 240-205222/27 | P4000027.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1221 Peak 1 | BNB | Lin1 | 1144216<br>32732245 | 1896892 | 3746401 | 8446168 | 16623076 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1221 Peak 2 | BNB | Lin1 | 779046<br>21219388 | 1287048 | 2589485 | 5697518 | 11015433 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1221 Peak 3 | BNB | Lin1 | 2852579<br>81639590 | 4640680 | 9222258 | 20939223 | 41609910 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 1 | BNB | Lin1 | 4010565<br>116813894 | 6484721 | 12487590 | 28845961 | 58294542 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 2 | BNB | Lin1 | 5987736<br>196460807 | 9760491 | 19563472 | 47143570 | 97436908 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 3 | BNB | Lin1 | 4572868<br>141075927 | 7400562 | 14247776 | 34764240 | 71829399 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 4 | BNB | Lin1 | 3496379<br>119217443 | 5617778 | 11273291 | 27783669 | 58818946 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 5 | BNB | Lin1 | 6220148<br>191684824 | 9984558 | 19492658 | 46853714 | 96578739 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4 GC Column: CLP-2 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 21:24 Calibration End Date: 11/04/2015 22:46 Calibration ID: 31286

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/28 | P4000028.D |
| Level 2 | STD01 240-205222/29 | P4000029.D |
| Level 3 | STD02 240-205222/30 | P4000030.D |
| Level 4 | STD05 240-205222/31 | P4000031.D |
| Level 5 | STD1 240-205222/32 | P4000032.D |
| Level 6 | STD2 240-205222/33 | P4000033.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1016 Peak 1 | 0.0178<br>0.0128 | 0.0153 | 0.0141 | 0.0136 | 0.0134 | Lin1 | 0.0003 | 0.0128 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 2 | 0.0346<br>0.0247 | 0.0296 | 0.0276 | 0.0265 | 0.0260 | Lin1 | 0.0005 | 0.0249 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1016 Peak 3 | 0.0686<br>0.0581 | 0.0603 | 0.0585 | 0.0592 | 0.0594 | Lin1 | 0.0004 | 0.0582 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 4 | 0.0299<br>0.0224 | 0.0253 | 0.0239 | 0.0233 | 0.0230 | Lin1 | 0.0004 | 0.0224 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 5 | 0.0240<br>0.0168 | 0.0200 | 0.0185 | 0.0179 | 0.0174 | Lin1 | 0.0004 | 0.0168 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 1 | 0.0436<br>0.0313 | 0.0346 | 0.0370 | 0.0366 | 0.0324 | Lin1 | 0.0006 | 0.0318 | | | | | | | 0.9970 | | 0.9900 |
| PCB-1260 Peak 2 | 0.0509<br>0.0407 | 0.0443 | 0.0434 | 0.0430 | 0.0419 | Lin1 | 0.0005 | 0.0409 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 3 | 0.0730<br>0.0619 | 0.0645 | 0.0639 | 0.0642 | 0.0633 | Lin1 | 0.0005 | 0.0622 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 4 | 0.0854<br>0.0747 | 0.0754 | 0.0759 | 0.0763 | 0.0752 | Lin1 | 0.0004 | 0.0747 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 5 | 0.0603<br>0.0505 | 0.0529 | 0.0527 | 0.0524 | 0.0514 | Lin1 | 0.0004 | 0.0506 | | | | | | | 1.0000 | | 0.9900 |
| Tetrachloro-m-xylene | 0.6148<br>0.6612 | 0.6230 | 0.6064 | 0.6560 | 0.6839 | Lin1 | 0 | 0.6687 | | | | | | | 0.9990 | | 0.9900 |
| DCB Decachlorobiphenyl | 0.6529<br>0.5587 | 0.6535 | 0.6238 | 0.6147 | 0.5968 | Lin1 | 0.0003 | 0.5721 | | | | | | | 0.9980 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4 GC Column: CLP-2 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 21:24 Calibration End Date: 11/04/2015 22:46 Calibration ID: 31286

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/28 | P4000028.D |
| Level 2 | STD01 240-205222/29 | P4000029.D |
| Level 3 | STD02 240-205222/30 | P4000030.D |
| Level 4 | STD05 240-205222/31 | P4000031.D |
| Level 5 | STD1 240-205222/32 | P4000032.D |
| Level 6 | STD2 240-205222/33 | P4000033.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1016 Peak 1 | BNB | Lin1 | 2505369<br>69882586 | 4111263 | 7796552 | 18785371 | 36011409 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 2 | BNB | Lin1 | 4860352<br>135341671 | 7952057 | 15208450 | 36697302 | 70055011 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 3 | BNB | Lin1 | 9633407<br>317864937 | 16196776 | 32276710 | 81841773 | 160099876 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 4 | BNB | Lin1 | 4196549<br>122500558 | 6792634 | 13186046 | 32174406 | 61997458 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 5 | BNB | Lin1 | 3366584<br>91951765 | 5377572 | 10199028 | 24741933 | 46920127 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 1 | BNB | Lin1 | 6117308<br>171297828 | 9282513 | 20424127 | 50590638 | 87425397 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 2 | BNB | Lin1 | 7143308<br>222482028 | 11912264 | 23947828 | 59394587 | 112955799 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 3 | BNB | Lin1 | 10249479<br>338705571 | 17336680 | 35280631 | 88729710 | 170599873 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 4 | BNB | Lin1 | 11993019<br>408685041 | 20262999 | 41909124 | 105539759 | 202783930 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 5 | BNB | Lin1 | 8470490<br>276063433 | 14204582 | 29098510 | 72460597 | 138608189 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| Tetrachloro-m-xylene | BNB | Lin1 | 4317714<br>180868998 | 8366942 | 16734439 | 45350885 | 92162505 | 0.00250<br>0.100 | 0.00500 | 0.0100 | 0.0250 | 0.0500 |
| DCB Decachlorobiphenyl | BNB | Lin1 | 4585127<br>152837892 | 8777594 | 17215039 | 42499125 | 80421640 | 0.00250<br>0.100 | 0.00500 | 0.0100 | 0.0250 | 0.0500 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-205222/39  Calibration Date: 11/05/2015 00:24

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 21:24

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 22:46

Lab File ID: P4000039.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0108 | | 0.822 | 1.00 | -17.8 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0217 | | 0.853 | 1.00 | -14.7 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0497 | | 0.848 | 1.00 | -15.2 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0191 | | 0.837 | 1.00 | -16.3 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0142 | | 0.824 | 1.00 | -17.6 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0258 | | 0.789 | 1.00 | -21.1* | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0324 | | 0.781 | 1.00 | -21.9* | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0430 | | 0.685 | 1.00 | -31.5* | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0708 | | 0.943 | 1.00 | -5.7 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0461 | | 0.903 | 1.00 | -9.7 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6638 | | 0.0500 | 0.0500 | -0.0 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.5091 | | 0.0439 | 0.0500 | -12.2 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-205222/39 Calibration Date: 11/05/2015 00:24

Instrument ID: A2HP4 Calib Start Date: 11/04/2015 21:24

GC Column: CLP-2 ID: 0.53(mm) Calib End Date: 11/04/2015 22:46

Lab File ID: P4000039.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1016 Peak 1 | 4.51 | 4.48 | 4.54 |
| PCB-1016 Peak 2 | 5.20 | 5.16 | 5.22 |
| PCB-1016 Peak 3 | 5.90 | 5.87 | 5.93 |
| PCB-1016 Peak 4 | 6.11 | 6.07 | 6.13 |
| PCB-1016 Peak 5 | 6.75 | 6.71 | 6.77 |
| PCB-1260 Peak 1 | 8.11 | 8.08 | 8.14 |
| PCB-1260 Peak 2 | 8.31 | 8.28 | 8.34 |
| PCB-1260 Peak 3 | 8.63 | 8.59 | 8.65 |
| PCB-1260 Peak 4 | 9.18 | 9.14 | 9.20 |
| PCB-1260 Peak 5 | 9.48 | 9.44 | 9.50 |
| Tetrachloro-m-xylene | 3.59 | 3.55 | 3.61 |
| DCB Decachlorobiphenyl | 10.41 | 10.38 | 10.44 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/3  Calibration Date: 11/11/2015 11:01

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 21:24

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 22:46

Lab File ID: P4000003.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0146 | | 0.549 | 0.500 | 9.8 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0282 | | 0.545 | 0.500 | 9.0 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0618 | | 0.525 | 0.500 | 5.0 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0246 | | 0.534 | 0.500 | 6.7 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0186 | | 0.530 | 0.500 | 6.1 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0333 | | 0.503 | 0.500 | 0.7 | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0427 | | 0.510 | 0.500 | 1.9 | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0632 | | 0.501 | 0.500 | 0.3 | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0741 | | 0.491 | 0.500 | -1.8 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0509 | | 0.494 | 0.500 | -1.2 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6986 | | 0.0264 | 0.0250 | 5.8 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.5950 | | 0.0254 | 0.0250 | 1.6 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/3 Calibration Date: 11/11/2015 11:01

Instrument ID: A2HP4 Calib Start Date: 11/04/2015 21:24

GC Column: CLP-2 ID: 0.53(mm) Calib End Date: 11/04/2015 22:46

Lab File ID: P4000003.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1016 Peak 1 | 4.49 | 4.46 | 4.52 |
| PCB-1016 Peak 2 | 5.18 | 5.15 | 5.21 |
| PCB-1016 Peak 3 | 5.88 | 5.85 | 5.91 |
| PCB-1016 Peak 4 | 6.09 | 6.06 | 6.12 |
| PCB-1016 Peak 5 | 6.73 | 6.70 | 6.76 |
| PCB-1260 Peak 1 | 8.09 | 8.06 | 8.12 |
| PCB-1260 Peak 2 | 8.30 | 8.27 | 8.33 |
| PCB-1260 Peak 3 | 8.61 | 8.58 | 8.64 |
| PCB-1260 Peak 4 | 9.16 | 9.13 | 9.19 |
| PCB-1260 Peak 5 | 9.46 | 9.43 | 9.49 |
| Tetrachloro-m-xylene | 3.57 | 3.54 | 3.60 |
| DCB Decachlorobiphenyl | 10.40 | 10.37 | 10.43 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/4  Calibration Date: 11/11/2015 11:17

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 14:51

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 16:13

Lab File ID: P4000004.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1232 Peak 1 | Lin1 | | 0.0149 | | 0.593 | 0.500 | 18.6 | 20.0 |
| PCB-1232 Peak 2 | Lin1 | | 0.0121 | | 0.604 | 0.500 | 20.8* | 20.0 |
| PCB-1232 Peak 3 | Lin1 | | 0.0235 | | 0.569 | 0.500 | 13.8 | 20.0 |
| PCB-1232 Peak 4 | Lin1 | | 0.0099 | | 0.592 | 0.500 | 18.4 | 20.0 |
| PCB-1232 Peak 5 | Lin1 | | 0.0080 | | 0.600 | 0.500 | 19.9 | 20.0 |
| PCB-1262 Peak 1 | Lin1 | | 0.0266 | | 0.539 | 0.500 | 7.8 | 20.0 |
| PCB-1262 Peak 2 | Lin1 | | 0.0407 | | 0.535 | 0.500 | 7.0 | 20.0 |
| PCB-1262 Peak 3 | Lin1 | | 0.0338 | | 0.527 | 0.500 | 5.5 | 20.0 |
| PCB-1262 Peak 4 | Lin1 | | 0.0787 | | 0.515 | 0.500 | 3.0 | 20.0 |
| PCB-1262 Peak 5 | Lin1 | | 0.0557 | | 0.517 | 0.500 | 3.3 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/4 Calibration Date: 11/11/2015 11:17

Instrument ID: A2HP4 Calib Start Date: 11/04/2015 14:51

GC Column: CLP-2 ID: 0.53(mm) Calib End Date: 11/04/2015 16:13

Lab File ID: P4000004.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1232 Peak 1 | 4.49 | 4.46 | 4.52 |
| PCB-1232 Peak 2 | 5.18 | 5.15 | 5.21 |
| PCB-1232 Peak 3 | 5.89 | 5.86 | 5.92 |
| PCB-1232 Peak 4 | 6.09 | 6.06 | 6.12 |
| PCB-1232 Peak 5 | 6.88 | 6.85 | 6.91 |
| PCB-1262 Peak 1 | 8.30 | 8.27 | 8.33 |
| PCB-1262 Peak 2 | 8.70 | 8.67 | 8.73 |
| PCB-1262 Peak 3 | 8.95 | 8.92 | 8.98 |
| PCB-1262 Peak 4 | 9.16 | 9.13 | 9.19 |
| PCB-1262 Peak 5 | 9.46 | 9.43 | 9.49 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/5  Calibration Date: 11/11/2015 11:34

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 16:30

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 17:51

Lab File ID: P4000005.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1242 Peak 1 | Lin1 | | 0.0100 | | 0.583 | 0.500 | 16.7 | 20.0 |
| PCB-1242 Peak 2 | Lin1 | | 0.0189 | | 0.571 | 0.500 | 14.1 | 20.0 |
| PCB-1242 Peak 3 | Lin1 | | 0.0402 | | 0.543 | 0.500 | 8.7 | 20.0 |
| PCB-1242 Peak 4 | Lin1 | | 0.0163 | | 0.557 | 0.500 | 11.4 | 20.0 |
| PCB-1242 Peak 5 | Lin1 | | 0.0127 | | 0.562 | 0.500 | 12.3 | 20.0 |
| PCB-1268 Peak 1 | Lin1 | | 0.0893 | | 0.490 | 0.500 | -1.9 | 20.0 |
| PCB-1268 Peak 2 | Lin1 | | 0.0866 | | 0.491 | 0.500 | -1.8 | 20.0 |
| PCB-1268 Peak 3 | Lin1 | | 0.0761 | | 0.496 | 0.500 | -0.8 | 20.0 |
| PCB-1268 Peak 4 | Lin1 | | 0.0306 | | 0.490 | 0.500 | -2.0 | 20.0 |
| PCB-1268 Peak 5 | Lin1 | | 0.2305 | | 0.468 | 0.500 | -6.4 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/5 Calibration Date: 11/11/2015 11:34

Instrument ID: A2HP4 Calib Start Date: 11/04/2015 16:30

GC Column: CLP-2 ID: 0.53(mm) Calib End Date: 11/04/2015 17:51

Lab File ID: P4000005.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1242 Peak 1 | 4.50 | 4.47 | 4.53 |
| PCB-1242 Peak 2 | 5.18 | 5.15 | 5.21 |
| PCB-1242 Peak 3 | 5.89 | 5.86 | 5.92 |
| PCB-1242 Peak 4 | 6.09 | 6.06 | 6.12 |
| PCB-1242 Peak 5 | 6.73 | 6.70 | 6.76 |
| PCB-1268 Peak 1 | 9.46 | 9.43 | 9.49 |
| PCB-1268 Peak 2 | 9.50 | 9.47 | 9.53 |
| PCB-1268 Peak 3 | 9.70 | 9.67 | 9.73 |
| PCB-1268 Peak 4 | 9.93 | 9.90 | 9.96 |
| PCB-1268 Peak 5 | 10.19 | 10.16 | 10.22 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/6  Calibration Date: 11/11/2015 11:50

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 18:08

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 19:30

Lab File ID: P4000006.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1248 Peak 1 | Lin1 | | 0.0255 | | 0.524 | 0.500 | 4.8 | 20.0 |
| PCB-1248 Peak 2 | Lin1 | | 0.0130 | | 0.533 | 0.500 | 6.6 | 20.0 |
| PCB-1248 Peak 3 | Lin1 | | 0.0193 | | 0.524 | 0.500 | 4.7 | 20.0 |
| PCB-1248 Peak 4 | Lin1 | | 0.0236 | | 0.516 | 0.500 | 3.2 | 20.0 |
| PCB-1248 Peak 5 | Lin1 | | 0.0119 | | 0.513 | 0.500 | 2.6 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/6 Calibration Date: 11/11/2015 11:50

Instrument ID: A2HP4 Calib Start Date: 11/04/2015 18:08

GC Column: CLP-2 ID: 0.53(mm) Calib End Date: 11/04/2015 19:30

Lab File ID: P4000006.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1248 Peak 1 | 5.88 | 5.85 | 5.91 |
| PCB-1248 Peak 2 | 6.44 | 6.41 | 6.47 |
| PCB-1248 Peak 3 | 6.73 | 6.70 | 6.76 |
| PCB-1248 Peak 4 | 7.23 | 7.20 | 7.26 |
| PCB-1248 Peak 5 | 7.91 | 7.88 | 7.94 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/7  Calibration Date: 11/11/2015 12:06

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 19:46

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 21:08

Lab File ID: P4000007.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1221 Peak 1 | Lin1 | | 0.0071 | | 0.575 | 0.500 | 15.0 | 20.0 |
| PCB-1221 Peak 2 | Lin1 | | 0.0048 | | 0.590 | 0.500 | 18.0 | 20.0 |
| PCB-1221 Peak 3 | Lin1 | | 0.0173 | | 0.565 | 0.500 | 13.1 | 20.0 |
| PCB-1254 Peak 1 | Lin1 | | 0.0240 | | 0.556 | 0.500 | 11.2 | 20.0 |
| PCB-1254 Peak 2 | Lin1 | | 0.0379 | | 0.530 | 0.500 | 6.1 | 20.0 |
| PCB-1254 Peak 3 | Lin1 | | 0.0280 | | 0.537 | 0.500 | 7.4 | 20.0 |
| PCB-1254 Peak 4 | Lin1 | | 0.0225 | | 0.526 | 0.500 | 5.1 | 20.0 |
| PCB-1254 Peak 5 | Lin1 | | 0.0372 | | 0.528 | 0.500 | 5.6 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/7 Calibration Date: 11/11/2015 12:06

Instrument ID: A2HP4 Calib Start Date: 11/04/2015 19:46

GC Column: CLP-2 ID: 0.53(mm) Calib End Date: 11/04/2015 21:08

Lab File ID: P4000007.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1221 Peak 1 | 4.15 | 4.12 | 4.18 |
| PCB-1221 Peak 2 | 4.39 | 4.36 | 4.42 |
| PCB-1221 Peak 3 | 4.50 | 4.47 | 4.53 |
| PCB-1254 Peak 1 | 7.23 | 7.20 | 7.26 |
| PCB-1254 Peak 2 | 7.91 | 7.88 | 7.94 |
| PCB-1254 Peak 3 | 8.13 | 8.10 | 8.16 |
| PCB-1254 Peak 4 | 8.45 | 8.42 | 8.48 |
| PCB-1254 Peak 5 | 8.62 | 8.59 | 8.65 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206283/28  Calibration Date: 11/11/2015 17:50

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 21:24

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 22:46

Lab File ID: P4000028.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0144 | | 0.541 | 0.500 | 8.2 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0279 | | 0.540 | 0.500 | 7.9 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0609 | | 0.517 | 0.500 | 3.4 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0243 | | 0.528 | 0.500 | 5.6 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0182 | | 0.519 | 0.500 | 3.9 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0327 | | 0.493 | 0.500 | -1.4 | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0419 | | 0.500 | 0.500 | 0.0 | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0623 | | 0.494 | 0.500 | -1.2 | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0728 | | 0.482 | 0.500 | -3.5 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0501 | | 0.486 | 0.500 | -2.7 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6986 | | 0.0264 | 0.0250 | 5.8 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.5154 | | 0.0219 | 0.0250 | -12.3 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206283/28 Calibration Date: 11/11/2015 17:50

Instrument ID: A2HP4 Calib Start Date: 11/04/2015 21:24

GC Column: CLP-2 ID: 0.53(mm) Calib End Date: 11/04/2015 22:46

Lab File ID: P4000028.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1016 Peak 1 | 4.51 | 4.48 | 4.54 |
| PCB-1016 Peak 2 | 5.19 | 5.16 | 5.22 |
| PCB-1016 Peak 3 | 5.90 | 5.87 | 5.93 |
| PCB-1016 Peak 4 | 6.10 | 6.07 | 6.13 |
| PCB-1016 Peak 5 | 6.74 | 6.71 | 6.77 |
| PCB-1260 Peak 1 | 8.11 | 8.08 | 8.14 |
| PCB-1260 Peak 2 | 8.31 | 8.28 | 8.34 |
| PCB-1260 Peak 3 | 8.62 | 8.59 | 8.65 |
| PCB-1260 Peak 4 | 9.17 | 9.14 | 9.20 |
| PCB-1260 Peak 5 | 9.47 | 9.44 | 9.50 |
| Tetrachloro-m-xylene | 3.58 | 3.55 | 3.61 |
| DCB Decachlorobiphenyl | 10.41 | 10.38 | 10.44 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205809/21-A

Matrix: Solid | Lab File ID: P4000026.D

Analysis Method: 8082A | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 17:17

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 12 | U | 33 | 12 |
| 11104-28-2 | Aroclor-1221 | 16 | U | 33 | 16 |
| 11141-16-5 | Aroclor-1232 | 20 | U | 33 | 20 |
| 53469-21-9 | Aroclor-1242 | 11 | U | 33 | 11 |
| 12672-29-6 | Aroclor-1248 | 8.0 | U | 33 | 8.0 |
| 11097-69-1 | Aroclor-1254 | 14 | U | 33 | 14 |
| 11096-82-5 | Aroclor-1260 | 9.0 | U | 33 | 9.0 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 75 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 89 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-205809/22-A

Matrix: Solid | Lab File ID: P4000027.D

Analysis Method: 8082A | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 17:33

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 196 | | 33 | 12 |
| 11096-82-5 | Aroclor-1260 | 201 | | 33 | 9.0 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 75 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 85 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: P4000030.D

Analysis Method: 8082A | Date Collected: 11/05/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.65(g) | Date Analyzed: 11/11/2015 18:22

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 244 | | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 260 | | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 81 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 95 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD | Lab Sample ID: 240-57597-15 MSD

Matrix: Solid | Lab File ID: P4000031.D

Analysis Method: 8082A | Date Collected: 11/05/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.78(g) | Date Analyzed: 11/11/2015 18:39

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 229 | | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.4 | U | 39 | 9.4 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 39 | 16 |
| 11096-82-5 | Aroclor-1260 | 238 | | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 77 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 89 | | 10-155 |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4 Start Date: 11/04/2015 14:51

Analysis Batch Number: 205222 End Date: 11/05/2015 00:57

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD005 240-205222/4 IC | | 11/04/2015 14:51 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/4 IC | | 11/04/2015 14:51 | 1 | P4000004.D | CLP-2 0.53(mm) |
| STD01 240-205222/5 IC | | 11/04/2015 15:08 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/5 IC | | 11/04/2015 15:08 | 1 | P4000005.D | CLP-2 0.53(mm) |
| STD02 240-205222/6 IC | | 11/04/2015 15:24 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/6 IC | | 11/04/2015 15:24 | 1 | P4000006.D | CLP-2 0.53(mm) |
| STD05 240-205222/7 IC | | 11/04/2015 15:40 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/7 IC | | 11/04/2015 15:40 | 1 | P4000007.D | CLP-2 0.53(mm) |
| STD1 240-205222/8 IC | | 11/04/2015 15:57 | 1 | | CLP-1 0.53(mm) |
| STD1 240-205222/8 IC | | 11/04/2015 15:57 | 1 | P4000008.D | CLP-2 0.53(mm) |
| STD2 240-205222/9 IC | | 11/04/2015 16:13 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/9 IC | | 11/04/2015 16:13 | 1 | P4000009.D | CLP-2 0.53(mm) |
| STD005 240-205222/10 IC | | 11/04/2015 16:30 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/10 IC | | 11/04/2015 16:30 | 1 | P4000010.D | CLP-2 0.53(mm) |
| STD01 240-205222/11 IC | | 11/04/2015 16:46 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/11 IC | | 11/04/2015 16:46 | 1 | P4000011.D | CLP-2 0.53(mm) |
| STD02 240-205222/12 IC | | 11/04/2015 17:02 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/12 IC | | 11/04/2015 17:02 | 1 | P4000012.D | CLP-2 0.53(mm) |
| STD05 240-205222/13 IC | | 11/04/2015 17:19 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/13 IC | | 11/04/2015 17:19 | 1 | P4000013.D | CLP-2 0.53(mm) |
| STD1 240-205222/14 IC | | 11/04/2015 17:35 | 1 | | CLP-1 0.53(mm) |
| STD1 240-205222/14 IC | | 11/04/2015 17:35 | 1 | P4000014.D | CLP-2 0.53(mm) |
| STD2 240-205222/15 IC | | 11/04/2015 17:51 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/15 IC | | 11/04/2015 17:51 | 1 | P4000015.D | CLP-2 0.53(mm) |
| STD005 240-205222/16 IC | | 11/04/2015 18:08 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/16 IC | | 11/04/2015 18:08 | 1 | P4000016.D | CLP-2 0.53(mm) |
| STD01 240-205222/17 IC | | 11/04/2015 18:24 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/17 IC | | 11/04/2015 18:24 | 1 | P4000017.D | CLP-2 0.53(mm) |
| STD02 240-205222/18 IC | | 11/04/2015 18:41 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/18 IC | | 11/04/2015 18:41 | 1 | P4000018.D | CLP-2 0.53(mm) |
| STD05 240-205222/19 IC | | 11/04/2015 18:57 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/19 IC | | 11/04/2015 18:57 | 1 | P4000019.D | CLP-2 0.53(mm) |
| STD1 240-205222/20 IC | | 11/04/2015 19:13 | 1 | | CLP-1 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4 Start Date: 11/04/2015 14:51

Analysis Batch Number: 205222 End Date: 11/05/2015 00:57

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-205222/20 IC | | 11/04/2015 19:13 | 1 | P4000020.D | CLP-2 0.53(mm) |
| STD2 240-205222/21 IC | | 11/04/2015 19:30 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/21 IC | | 11/04/2015 19:30 | 1 | P4000021.D | CLP-2 0.53(mm) |
| STD005 240-205222/22 IC | | 11/04/2015 19:46 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/22 IC | | 11/04/2015 19:46 | 1 | P4000022.D | CLP-2 0.53(mm) |
| STD01 240-205222/23 IC | | 11/04/2015 20:02 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/23 IC | | 11/04/2015 20:02 | 1 | P4000023.D | CLP-2 0.53(mm) |
| STD02 240-205222/24 IC | | 11/04/2015 20:19 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/24 IC | | 11/04/2015 20:19 | 1 | P4000024.D | CLP-2 0.53(mm) |
| STD05 240-205222/25 IC | | 11/04/2015 20:35 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/25 IC | | 11/04/2015 20:35 | 1 | P4000025.D | CLP-2 0.53(mm) |
| STD1 240-205222/26 IC | | 11/04/2015 20:51 | 1 | | CLP-1 0.53(mm) |
| STD1 240-205222/26 IC | | 11/04/2015 20:51 | 1 | P4000026.D | CLP-2 0.53(mm) |
| STD2 240-205222/27 IC | | 11/04/2015 21:08 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/27 IC | | 11/04/2015 21:08 | 1 | P4000027.D | CLP-2 0.53(mm) |
| STD005 240-205222/28 IC | | 11/04/2015 21:24 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/28 IC | | 11/04/2015 21:24 | 1 | P4000028.D | CLP-2 0.53(mm) |
| STD01 240-205222/29 IC | | 11/04/2015 21:40 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/29 IC | | 11/04/2015 21:40 | 1 | P4000029.D | CLP-2 0.53(mm) |
| STD02 240-205222/30 IC | | 11/04/2015 21:57 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/30 IC | | 11/04/2015 21:57 | 1 | P4000030.D | CLP-2 0.53(mm) |
| STD05 240-205222/31 ICIS | | 11/04/2015 22:13 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/31 ICIS | | 11/04/2015 22:13 | 1 | P4000031.D | CLP-2 0.53(mm) |
| STD1 240-205222/32 IC | | 11/04/2015 22:29 | 1 | | CLP-1 0.53(mm) |
| STD1 240-205222/32 IC | | 11/04/2015 22:29 | 1 | P4000032.D | CLP-2 0.53(mm) |
| STD2 240-205222/33 IC | | 11/04/2015 22:46 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/33 IC | | 11/04/2015 22:46 | 1 | P4000033.D | CLP-2 0.53(mm) |
| ICV 240-205222/34 | | 11/04/2015 23:02 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/34 | | 11/04/2015 23:02 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/35 | | 11/04/2015 23:19 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/35 | | 11/04/2015 23:19 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/36 | | 11/04/2015 23:35 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/36 | | 11/04/2015 23:35 | 1 | | CLP-2 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4  Start Date: 11/04/2015 14:51

Analysis Batch Number: 205222  End Date: 11/05/2015 00:57

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| ICV 240-205222/37 | | 11/04/2015 23:51 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/37 | | 11/04/2015 23:51 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/38 | | 11/05/2015 00:08 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/38 | | 11/05/2015 00:08 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/39 | | 11/05/2015 00:24 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/39 | | 11/05/2015 00:24 | 1 | P4000039.D | CLP-2 0.53(mm) |
| ICV 240-205222/40 | | 11/05/2015 00:40 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/40 | | 11/05/2015 00:40 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/41 | | 11/05/2015 00:57 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/41 | | 11/05/2015 00:57 | 1 | | CLP-2 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4 Start Date: 11/11/2015 11:01

Analysis Batch Number: 206283 End Date: 11/11/2015 18:39

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-206283/3 CCVIS | | 11/11/2015 11:01 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/3 CCVIS | | 11/11/2015 11:01 | 1 | P4000003.D | CLP-2 0.53(mm) |
| CCV 240-206283/4 | | 11/11/2015 11:17 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/4 | | 11/11/2015 11:17 | 1 | P4000004.D | CLP-2 0.53(mm) |
| CCV 240-206283/5 | | 11/11/2015 11:34 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/5 | | 11/11/2015 11:34 | 1 | P4000005.D | CLP-2 0.53(mm) |
| CCV 240-206283/6 | | 11/11/2015 11:50 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/6 | | 11/11/2015 11:50 | 1 | P4000006.D | CLP-2 0.53(mm) |
| CCV 240-206283/7 | | 11/11/2015 12:06 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/7 | | 11/11/2015 12:06 | 1 | P4000007.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 13:28 | 1 | | CLP-1 0.53(mm) |
| 240-57597-1 | F-VB-UST-1-14 | 11/11/2015 13:28 | 1 | P4000012.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 13:45 | 1 | | CLP-1 0.53(mm) |
| 240-57597-2 | F-VB-UST-2-14 | 11/11/2015 13:45 | 1 | P4000013.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 14:01 | 1 | | CLP-1 0.53(mm) |
| 240-57597-3 | F-VB-UST-3-14 | 11/11/2015 14:01 | 1 | P4000014.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 14:17 | 1 | | CLP-1 0.53(mm) |
| 240-57597-4 | F-VB-UST-4-14 | 11/11/2015 14:17 | 1 | P4000015.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 14:33 | 1 | | CLP-1 0.53(mm) |
| 240-57597-5 | F-VB-UST-5-14 | 11/11/2015 14:33 | 1 | P4000016.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 14:50 | 1 | | CLP-1 0.53(mm) |
| 240-57597-6 | F-VB-UST-6-14 | 11/11/2015 14:50 | 1 | P4000017.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 15:06 | 1 | | CLP-1 0.53(mm) |
| 240-57597-7 | F-VB-UST-7-14 | 11/11/2015 15:06 | 1 | P4000018.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 15:23 | 1 | | CLP-1 0.53(mm) |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 15:23 | 1 | P4000019.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 15:39 | 1 | | CLP-1 0.53(mm) |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 15:39 | 1 | P4000020.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 15:55 | 1 | | CLP-1 0.53(mm) |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 15:55 | 1 | P4000021.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 16:12 | 1 | | CLP-1 0.53(mm) |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 16:12 | 1 | P4000022.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 16:28 | 1 | | CLP-1 0.53(mm) |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 16:28 | 1 | P4000023.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 16:44 | 1 | | CLP-1 0.53(mm) |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 16:44 | 1 | P4000024.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 17:01 | 1 | | CLP-1 0.53(mm) |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 17:01 | 1 | P4000025.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 17:17 | 1 | | CLP-1 0.53(mm) |
| MB 240-205809/21-A | | 11/11/2015 17:17 | 1 | P4000026.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 17:33 | 1 | | CLP-1 0.53(mm) |
| LCS 240-205809/22-A | | 11/11/2015 17:33 | 1 | P4000027.D | CLP-2 0.53(mm) |
| CCVIS 240-206283/28 | | 11/11/2015 17:50 | 1 | | CLP-1 0.53(mm) |
| CCVIS 240-206283/28 | | 11/11/2015 17:50 | 1 | P4000028.D | CLP-2 0.53(mm) |

8082A

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4 Start Date: 11/11/2015 11:01

Analysis Batch Number: 206283 End Date: 11/11/2015 18:39

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| ZZZZZ | | 11/11/2015 18:06 | 1 | | CLP-1 0.53(mm) |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 18:06 | 1 | P4000029.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 18:22 | 1 | | CLP-1 0.53(mm) |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 18:22 | 1 | P4000030.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 18:39 | 1 | | CLP-1 0.53(mm) |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 18:39 | 1 | P4000031.D | CLP-2 0.53(mm) |

# PCBS BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205809  Batch Start Date: 11/09/15 11:00  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/10/15 03:00

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | ex10PPMSPK 00018 | ex2/.2SURRW 00062 | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-E-1 | F-VB-UST-1-14 | 3540C, 8082A | T | 29.54 g | 10 mL | | 1 mL | | |
| 240-57597-D-2 | F-VB-UST-2-14 | 3540C, 8082A | T | 30.07 g | 10 mL | | 1 mL | | |
| 240-57597-E-3 | F-VB-UST-3-14 | 3540C, 8082A | T | 30.14 g | 10 mL | | 1 mL | | |
| 240-57597-E-4 | F-VB-UST-4-14 | 3540C, 8082A | T | 29.54 g | 10 mL | | 1 mL | | |
| 240-57597-E-5 | F-VB-UST-5-14 | 3540C, 8082A | T | 30.48 g | 10 mL | | 1 mL | | |
| 240-57597-E-6 | F-VB-UST-6-14 | 3540C, 8082A | T | 29.87 g | 10 mL | | 1 mL | | |
| 240-57597-E-7 | F-VB-UST-7-14 | 3540C, 8082A | T | 30.04 g | 10 mL | | 1 mL | | |
| 240-57597-D-8 | F-VB-UST-8-14 | 3540C, 8082A | T | 29.74 g | 10 mL | | 1 mL | | |
| 240-57597-E-9 | F-VB-UST-9-14 | 3540C, 8082A | T | 30.07 g | 10 mL | | 1 mL | | |
| 240-57597-E-10 | F-VB-UST-10-14 | 3540C, 8082A | T | 29.97 g | 10 mL | | 1 mL | | |
| 240-57597-D-11 | F-VB-UST-11-14 | 3540C, 8082A | T | 30.28 g | 10 mL | | 1 mL | | |
| 240-57597-D-12 | F-VB-UST-12-14 | 3540C, 8082A | T | 30.29 g | 10 mL | | 1 mL | | |
| 240-57597-D-13 | F-VB-UST-13-14 | 3540C, 8082A | T | 30.41 g | 10 mL | | 1 mL | | |
| 240-57597-E-14 | F-VB-UST-14-14 | 3540C, 8082A | T | 30.19 g | 10 mL | | 1 mL | | |
| 240-57597-E-15 | F-VB-UST-DUP | 3540C, 8082A | T | 30.34 g | 10 mL | | 1 mL | | |
| 240-57597-E-15 MS | F-VB-UST-DUP | 3540C, 8082A | T | 29.65 g | 10 mL | 1 mL | 1 mL | | |
| 240-57597-E-15 MSD | F-VB-UST-DUP | 3540C, 8082A | T | 29.78 g | 10 mL | 1 mL | 1 mL | | |
| MB 240-205809/21 | | 3540C, 8082A | | 30 g | 10 mL | | 1 mL | | |
| LCS 240-205809/22 | | 3540C, 8082A | | 30 g | 10 mL | 1 mL | 1 mL | | |

| Batch Notes | |
|---|---|
| Acid used for Clean Up Reagent | 2226515 |
| Balance ID | 1339912 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2247000 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GENERAL CHEMISTRY

COVER PAGE
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job Number: 240-57597-1

SDG No.:

Project: MRC Soil Remedy Block F UST

| Client Sample ID | Lab Sample ID |
|---|---|
| F-VB-UST-1-14 | 240-57597-1 |
| F-VB-UST-2-14 | 240-57597-2 |
| F-VB-UST-3-14 | 240-57597-3 |
| F-VB-UST-4-14 | 240-57597-4 |
| F-VB-UST-5-14 | 240-57597-5 |
| F-VB-UST-6-14 | 240-57597-6 |
| F-VB-UST-7-14 | 240-57597-7 |
| F-VB-UST-8-14 | 240-57597-8 |
| F-VB-UST-9-14 | 240-57597-9 |
| F-VB-UST-10-14 | 240-57597-10 |
| F-VB-UST-11-14 | 240-57597-11 |
| F-VB-UST-12-14 | 240-57597-12 |
| F-VB-UST-13-14 | 240-57597-13 |
| F-VB-UST-14-14 | 240-57597-14 |
| F-VB-UST-DUP | 240-57597-15 |

Comments:

9-IN
DETECTION LIMITS
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job Number: 240-57597-1

SDG Number:

Matrix: Solid Instrument ID: NOEQUIP

Method: Moisture RL Date: 01/28/2010 09:24

| Analyte | Wavelength/ Mass | RL (%) | |
|---|---|---|---|
| Percent Moisture | | 0.1 | |
| Percent Solids | | 0.1 | |

9-IN
CALIBRATION BLANK DETECTION LIMITS
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job Number: 240-57597-1

SDG Number:

Matrix: Solid Instrument ID: NOEQUIP

Method: Moisture XRL Date: 01/28/2010 09:24

| Analyte | Wavelength/ Mass | XRL (mg/L) | |
|---|---|---|---|
| Percent Moisture | | 10 | |
| Percent Solids | | 10 | |

13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: NOEQUIP Method: Moisture

Start Date: 11/09/2015 10:19 End Date: 11/09/2015 10:19

| Lab Sample ID | D / F | Type | Time | Analytes | | | | | | | | | | | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | | % Sol | Moist | | | | | | | | | | | | | | | | | | |
| 240-57597-1 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-1 DU | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-2 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-3 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-4 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-5 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-6 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-7 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-8 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-9 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-10 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-10 DU | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-11 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-12 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-13 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-14 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
| 240-57597-15 | 1 | T | 10:19 | X | X | | | | | | | | | | | | | | | | | | |
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13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: NOEQUIP Method: Moisture

Start Date: 11/09/2015 10:19 End Date: 11/09/2015 10:19

| Lab Sample ID | D / F | Type | Time | Analytes | | | | | | | | | | | | | | | | | | | | |
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| | | | | % Sol | Moist | | | | | | | | | | | | | | | | | | | |
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13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: NOEQUIP Method: Moisture

Start Date: 11/09/2015 10:19 End Date: 11/09/2015 10:19

| Lab Sample ID | D / F | Type | Time | Analytes | | | | | | | | | | | | | | | | | | | | |
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| | | | | % Sol | Moist | | | | | | | | | | | | | | | | | | | |
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| ZZZZZZ | | | 10:19 | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 10:19 | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 10:19 | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 10:19 | | | | | | | | | | | | | | | | | | | | | |

Prep Types

T = Total/NA

# GENERAL CHEMISTRY BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205828  Batch Start Date: 11/09/15 10:19  Batch Analyst: Nolle, Laura C

Batch Method: Moisture  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | DishWeight | SampleMassWet | SampleMassDry | | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-A-1 | F-VB-UST-1-14 | Moisture | T | 4.3999 g | 10.3790 g | 9.4392 g | | | |
| 240-57597-A-1 DU | F-VB-UST-1-14 | Moisture | T | 4.3999 g | 7.3944 g | 6.9197 g | | | |
| 240-57597-A-2 | F-VB-UST-2-14 | Moisture | T | 4.3999 g | 15.6278 g | 13.9472 g | | | |
| 240-57597-A-3 | F-VB-UST-3-14 | Moisture | T | 4.3999 g | 18.5326 g | 16.0226 g | | | |
| 240-57597-A-4 | F-VB-UST-4-14 | Moisture | T | 4.3999 g | 12.1776 g | 10.9626 g | | | |
| 240-57597-A-5 | F-VB-UST-5-14 | Moisture | T | 4.3999 g | 11.5850 g | 10.4924 g | | | |
| 240-57597-A-6 | F-VB-UST-6-14 | Moisture | T | 4.3999 g | 16.0347 g | 14.1108 g | | | |
| 240-57597-A-7 | F-VB-UST-7-14 | Moisture | T | 4.3999 g | 15.3864 g | 13.6038 g | | | |
| 240-57597-A-8 | F-VB-UST-8-14 | Moisture | T | 4.3999 g | 12.8669 g | 11.3431 g | | | |
| 240-57597-A-9 | F-VB-UST-9-14 | Moisture | T | 4.3999 g | 14.6966 g | 12.7294 g | | | |
| 240-57597-A-10 | F-VB-UST-10-14 | Moisture | T | 4.3999 g | 12.7867 g | 11.6604 g | | | |
| 240-57597-A-10 DU | F-VB-UST-10-14 | Moisture | T | 4.3999 g | 8.0824 g | 7.5869 g | | | |
| 240-57597-A-11 | F-VB-UST-11-14 | Moisture | T | 4.3999 g | 14.0661 g | 12.4699 g | | | |
| 240-57597-A-12 | F-VB-UST-12-14 | Moisture | T | 4.3999 g | 14.4309 g | 12.8362 g | | | |
| 240-57597-A-13 | F-VB-UST-13-14 | Moisture | T | 4.3999 g | 11.8606 g | 10.7651 g | | | |
| 240-57597-A-14 | F-VB-UST-14-14 | Moisture | T | 4.3999 g | 12.8355 g | 11.6131 g | | | |
| 240-57597-A-15 | F-VB-UST-DUP | Moisture | T | 4.3999 g | 14.0784 g | 12.6776 g | | | |

| Batch Notes | |
|---|---|
| Balance ID | B047 No Unit |
| Date and Time Samples in Desiccator | 0600 11/9/15 |
| Date and Time Samples out of Desiccator | 11/9/15 0930 |
| Date samples were placed in the oven | 11/6/15 |
| Oven Temp when samples are put in oven | 104 Degrees C |
| Date samples were removed from oven | 11/9/15 |
| Oven Temp when samples removed from oven | 104 Degrees C |
| Time Samples were removed from oven | 0600 |
| Oven ID | 002 |
| ID number of the thermometer | TempGuard Box-C #6 |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GENERAL CHEMISTRY BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205828  Batch Start Date: 11/09/15 10:19  Batch Analyst: Nolle, Laura C

Batch Method: Moisture  Batch End Date:

| Basis | Basis Description |
|---|---|
| T | Total/NA |

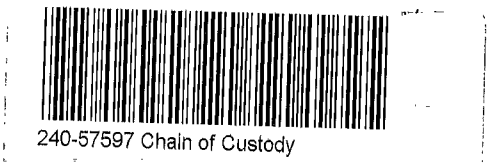
The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Shipping and Receiving Documents

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

TestAmerica Laboratories, Inc.

# CHAIN OF CUSTODY
# AND
# RECEIVING DOCUMENTS





4101 Shuffel Street, N.W. North Canton, OH 44720 **tel** 330.497.9396 **fax** 330.497.0772 **www.testamericainc.com**

0.8/C0.9
2.1/C2.2

# Chain of Custody Record #46 016125

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING
TestAmerica Laboratories, Inc.
TAL-8210 (0713)

Regulatory Program: ☐ DW ☐ NPDES ☐ RCRA ☐ Other:

| Client Contact | Project Manager: SCOTT NESBIT | Site Contact: J. MULLIS | Date: 11/5/15 | COC No: |
|---|---|---|---|---|
| Company Name: TETRA TECH | Tel/Fax: 412-559-8133 | Lab Contact: J. MCFADDEN | Carrier: FedEx | 1 of 2 COCs |
| Address: 20251 CENTURY BLVD STE 200 | Analysis Turnaround Time | | | Sampler: |
| City/State/Zip: GERMANTOWN, MD, 20874 | ☐ CALENDAR DAYS ☐ WORKING DAYS | | | For Lab Use Only: |
| Phone: 301-528-3000 | TAT if different from Below _____ | | | Walk-in Client: |
| Fax: 301-528-3000 | ☐ 2 weeks | | | Lab Sampling: |
| Project Name: MRC SOIL REMENT | ☒ 1 week | | | |
| Site: BLOCK F UST | ☐ 2 days | | | Job / SDG No.: |
| PO # 112IC07336 | ☐ 1 day | | | |

| Sample Identification | Sample Date | Sample Time | Sample Type (C=Comp, G=Grab) | Matrix | # of Cont. | Filtered Sample (Y/N) | Perform MS / MSD (Y/N) | VOCs (8260C) | SVOCs (8270D) | TPH GRO/DRO (8015B) | TCLP Metals (1311 then 6010) | PCBs (8082A) | Ignitability | Sample Specific Notes: |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| F-VB-UST-1-14 | 11/5/15 | 0900 | G | SO | 4 | | | X | X | X | X | X | | |
| F-VB-UST-2-14 | | 0902 | | | | | | X | X | X | X | X | | |
| F-VB-UST-3-14 | | 0904 | | | | | | X | X | X | X | X | | |
| F-VB-UST-4-14 | | 0906 | | | | | | X | X | X | X | X | | |
| F-VB-UST-5-14 | | 0908 | | | | | | X | X | X | X | X | | |
| F-VB-UST-6-14 | | 0910 | | | | | | X | X | X | X | X | | |
| F-VB-UST-7-14 | | 0912 | | | | | | X | X | X | X | X | | |
| F-VB-UST-8-14 | | 0914 | | | | | | X | X | X | X | X | | |
| F-VB-UST-9-14 | | 0916 | | | | | | X | X | X | X | X | | |
| F-VB-UST-10-14 | | 0918 | | | | | | X | X | X | X | X | | |
| F-VB-UST-11-14 | | 0920 | | | | | | X | X | X | X | X | | |
| F-VB-UST-12-14 | | 0922 | | | | | | X | X | X | X | X | | |

Preservation Used: 1= Ice, 2= HCl; 3= H2SO4; 4=HNO3; 5=NaOH; 6= Other _____

**Possible Hazard Identification:**
Are any samples from a listed EPA Hazardous Waste? Please List any EPA Waste Codes for the sample in the Comments Section if the lab is to dispose of the sample.

☒ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

**Sample Disposal ( A fee may be assessed if samples are retained longer than 1 month)**

☐ Return to Client ☒ Disposal by Lab ☐ Archive for _____ Months

**Special Instructions/QC Requirements & Comments:** MRC SOIL REMENT: BLOCK F UST BASE CONFIRMATION SAMPLES

| Custody Seals Intact: ☐ Yes ☐ No | Custody Seal No.: | | Cooler Temp. (°C): Obs'd: _____ Corr'd: _____ | | Therm ID No.: _____ |
|---|---|---|---|---|---|
| Relinquished by: | Company: TETRA TECH | Date/Time: 11/5/15 1530 | Received by: | Company: TestAmerica | Date/Time: 11/5/15 1530 |
| Relinquished by: | Company: TestAmerica | Date/Time: 11/5/15 1618 | Received by: | Company: TA | Date/Time: 11.6.15 1020 |
| Relinquished by: | Company: | Date/Time: | Received in Laboratory by: | Company: | Date/Time: |

# Chain of Custody Record

#46

016124

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING
TestAmerica Laboratories, Inc.
TAL-8210 (0713)

Regulatory Program: ☐ DW ☐ NPDES ☐ RCRA ☐ Other:

| Client Contact | Project Manager: SCOTT NESBIT | Site Contact: J. Mullis | Date: 11/5/15 | COC No: |
|---|---|---|---|---|
| Company Name: TETRA TECH | Tel/Fax: 412-559-8133 | Lab Contact: J. McFadden | Carrier: FedEx | 2 of 2 COCs |
| Address: 20251 Century Blvd. Ste. 200 | Analysis Turnaround Time | | | Sampler: |
| City/State/Zip: Germantown, MD, 20874 | ☐ CALENDAR DAYS ☐ WORKING DAYS | | | For Lab Use Only: |
| Phone: 301-528-3000 | TAT if different from Below _____ | | | Walk-in Client: |
| Fax: 301-528-3000 | ☐ 2 weeks | | | Lab Sampling: |
| Project Name: MRC SOIL REMEDY | ☒ 1 week | | | |
| Site: BLOCK F UST CONFIRMATION | ☐ 2 days | | | Job / SDG No.: |
| P O # 112IC07336 | ☐ 1 day | | | |

| Sample Identification | Sample Date | Sample Time | Sample Type (C=Comp, G=Grab) | Matrix | # of Cont. | Filtered Sample (Y/N) | Perform MS / MSD (Y/N) | VOCs (8260C) | SVOCs (8270D) | TPH GRO/DRO (8015B) | TCLP Metals (1311+6010) | RCRA (8015B) | Sample Specific Notes: |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| F-VB-UST-13-14 | 11/5/15 | 0924 | G | SO | 4 | | | X | X | X | X | X | |
| F-VB-UST-14-14 | | 0926 | | | | | | X | X | X | X | X | |
| F-VB-UST-DUP | | — | | | | | | X | X | X | X | X | # DUPLICATE |
| TB-110515 | | — | | AQ | 2 | | | X | | | | | |

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11/5/15

Preservation Used: (1= Ice) 2= HCl; 3= H2SO4; 4=HNO3; 5=NaOH; 6= Other

**Possible Hazard Identification:**
Are any samples from a listed EPA Hazardous Waste? Please List any EPA Waste Codes for the sample in the Comments Section if the lab is to dispose of the sample.

☒ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

**Sample Disposal ( A fee may be assessed if samples are retained longer than 1 month)**

☐ Return to Client ☒ Disposal by Lab ☐ Archive for _____ Months

**Special Instructions/QC Requirements & Comments:** SEE PAGE 1

| Custody Seals Intact: ☐ Yes ☐ No | Custody Seal No.: | | Cooler Temp. (°C): Obs'd: _____ Corr'd: _____ Therm ID No.: _____ | | |
|---|---|---|---|---|---|
| Relinquished by: | Company: TETRA TECH | Date/Time: 11/5/15 1530 | Received by: | Company: TestAmerica | Date/Time: 11/5/15 1530 |
| Relinquished by: | Company: TestAmerica | Date/Time: 11/5/15 1618 | Received by: | Company: TA | Date/Time: 11-6-15 1020 |
| Relinquished by: | Company: | Date/Time: | Received in Laboratory by: | Company: | Date/Time: |

**TestAmerica Canton Sample Receipt Form/Narrative** Login # : 57507
**Canton Facility**

Client Tetra Tech Site Name _____ Cooler unpacked by:

Cooler Received on 11-6-15 Opened on 11-6-15

FedEx: 1st Grd ~~Exp~~ UPS FAS Stetson Client Drop Off TestAmerica Courier Other _____

**Receipt After-hours**: Drop-off Date/Time _____ Storage Location _____

TestAmerica Cooler # _____ Foam Box Client Cooler Box ~~Other~~ _____

Packing material used: ~~Bubble Wrap~~ Foam Plastic Bag None Other _____

COOLANT: ~~Wet Ice~~ Blue Ice Dry Ice Water None

1. Cooler temperature upon receipt
   IR GUN# 53 (CF **+0.1** °C) Observed Cooler Temp._____°C Corrected Cooler Temp._____°C
   IR GUN# 48 (CF **-0.3** °C) Observed Cooler Temp._____°C Corrected Cooler Temp._____°C ☑ See Multiple Cooler Form
   ~~IR GUN# 5 (CF +0.4 °C) Observed Cooler Temp._____°C Corrected Cooler Temp._____°C~~
   IR GUN# 8 (CF **-0.5** °C) Observed Cooler Temp._____°C Corrected Cooler Temp._____°C

| | | |
|---|---|---|
| 2. Were custody seals on the outside of the cooler(s)? If Yes Quantity 2 | ~~Yes~~ | No |
| -Were custody seals on the outside of the cooler(s) signed & dated? | ~~Yes~~ | No NA |
| -Were custody seals on the bottle(s) or bottle kits (LLHg/MeHg)? | Yes | ~~No~~ |
| 3. Shippers' packing slip attached to the cooler(s)? | ~~Yes~~ | No |
| 4. Did custody papers accompany the sample(s)? | ~~Yes~~ | No |
| 5. Were the custody papers relinquished & signed in the appropriate place? | ~~Yes~~ | No |
| 6. Was/were the person(s) who collected the samples clearly identified on the COC? | Yes | ~~No~~ |
| 7. Did all bottles arrive in good condition (Unbroken)? | ~~Yes~~ | No |
| 8. Could all bottle labels be reconciled with the COC? | ~~Yes~~ | No |
| 9. Were correct bottle(s) used for the test(s) indicated? | ~~Yes~~ | No |
| 10. Sufficient quantity received to perform indicated analyses? | ~~Yes~~ | No |
| 11. Were sample(s) at the correct pH upon receipt? | Yes | No ~~NA~~ pH Strip Lot# **HC554612** |
| 12. Were VOAs on the COC? | ~~Yes~~ | No |
| 13. Were air bubbles >6 mm in any VOA vials? | Yes | ~~No~~ NA |
| 14. Was a trip blank present in the cooler(s)? Trip Blank Lot #_____ | ~~Yes~~ | No |

Contacted PM _____ Date _____ by _____ via Verbal Voice Mail Other

Concerning _____

**14. CHAIN OF CUSTODY & SAMPLE DISCREPANCIES** Samples processed by:

**15. SAMPLE CONDITION**

Sample(s) _____ were received after the recommended holding time had expired.

Sample(s) _____ were received in a broken container.

Sample(s) _____ were received with bubble >6 mm in diameter. (Notify PM)

**16. SAMPLE PRESERVATION**

Sample(s) _____ were further preserved in the laboratory.

Time preserved: _____ Preservative(s) added/Lot number(s): _____

TestAmerica Multiple Cooler Receipt Form/Narrative
Canton Facility

Login # : 575097

| Cooler # | IR Gun # | Observed Temp °C | Corrected Temp °C | Coolant |
|---|---|---|---|---|
| Client | 53 | 0.8 | 0.9 | ice |
| 1 | 1 | 2.1 | 2.2 | 1 |
| | | | | |
| | | | | |
| | | | |

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

# ANALYTICAL REPORT

Job Number: 240-57930-1

Job Description: LM MRC Block F UST Pipe Inv.

For:
Tetra Tech, Inc.
Foster Plaza 7
661 Anderson Drive
Pittsburgh, PA 15220-2745

Attention: Mr. Scott Nesbit

Approved for release.
John McFadden
Project Manager I
11/30/2015 1:57 PM

John McFadden, Project Manager I
4101 Shuffel Street NW, North Canton, OH, 44720
john.mcfadden@testamericainc.com
11/30/2015
Revision: 1

cc: Tony Apanavage
Samantha Brenner
Kelly Carper
Tobrena Sedlmyer

The test results in this report meet all NELAP requirements for parameters for which accreditation is required or available. Any exceptions to the NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory. This report is confidential and is intended for the sole use of TestAmerica and its client. All questions regarding this report should be directed to the TestAmerica Project Manager who has signed this report.

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# CASE NARRATIVE

## Client: Tetra Tech, Inc.

## Project: LM MRC Block F UST Pipe Inv.

## Report Number: 240-57930-1
## Revision 1

This report was revised on 11/30/15 to include the Form III's in the final report that were initially not included due to a linking issue with the laboratory's computer system.

With the exceptions noted as flags or footnotes, standard analytical protocols were followed in the analysis of the samples and no problems were encountered or anomalies observed. In addition all laboratory quality control samples were within established control limits, with any exceptions noted below. Each sample was analyzed to achieve the lowest possible reporting limit within the constraints of the method. In some cases, due to interference or analytes present at high concentrations, samples were diluted. For diluted samples, the reporting limits are adjusted relative to the dilution required.

TestAmerica Canton attests to the validity of the laboratory data generated by TestAmerica facilities reported herein. All analyses performed by TestAmerica facilities were done using established laboratory SOPs that incorporate QA/QC procedures described in the application methods. TestAmerica's operations groups have reviewed the data for compliance with the laboratory QA/QC plan, and data have been found to be compliant with laboratory protocols unless otherwise noted below.

The test results in this report meet all NELAP requirements for parameters for which accreditation is required or available. Any exceptions to NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory.

Calculations are performed before rounding to avoid round-off errors in calculated results.

All holding times were met and proper preservation noted for the methods performed on these samples, unless otherwise detailed in the individual sections below.

All solid sample results are reported on an "as received" basis unless otherwise indicated by the presence of a % solids value in the method header.

This laboratory report is confidential and is intended for the sole use of TestAmerica and its client.

**RECEIPT**

The samples were received on 11/14/2015 9:45 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 3.4º C.

**VOLATILE ORGANIC COMPOUNDS (GCMS)**

Samples F-VS-UST-P-1 (240-57930-1), F-VS-UST-P-2 (240-57930-2), F-VS-UST-P-3 (240-57930-3), F-VS-UST-P-4 (240-57930-4), F-VB-UST-P-1 (240-57930-5), F-VB-UST-P-2 (240-57930-6), F-VB-UST-P-3 (240-57930-7), F-VB-UST-P-4 (240-57930-8), F-VB-UST-P-5 (240-57930-9), F-VB-UST-P-6 (240-57930-10) and F-VS-UST-P-DUP (240-57930-11) were analyzed for volatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8260C. The samples were prepared and analyzed on 11/18/2015, 11/19/2015 and 11/20/2015.

Toluene was detected in method blanks MB 240-207343/6 and MB 240-207560/6 at a level that was above the method detection limit but below the reporting limit. The value should be considered an estimate, and has been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

Acetone was detected in method blank MB 240-207720/8 at a level exceeding the reporting limit. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

The continuing calibration verification (CCV) analyzed in batch 240-207343 was outside the method criteria for the following analytes: Carbon Tetrachloride, 2-Hexanone, 2-Methyl-2-propanol, 2-Chloroethyl Vinyl Ether, 4-Methyl-2-pentanone, 1,1,1-Trichloroethane, 2,2-Dichloropropane, Trichlorofluoromethane, 2-Butanone, and tert-Amyl Methyl Ether. An MRL standard at or below the reporting limit (RL) was analyzed with the affected samples: F-VS-UST-P-1 (240-57930-1), F-VS-UST-P-3 (240-57930-3), F-VB-UST-P-1 (240-57930-5), F-VB-UST-P-2 (240-57930-6), F-VB-UST-P-3 (240-57930-7), F-VB-UST-P-4 (240-57930-8), F-VB-UST-P-5 (240-57930-9), F-VB-UST-P-6 (240-57930-10), F-VS-UST-P-DUP (240-57930-11), (CCV 240-207343/4), (CCVIS 240-207343/3), (LCS 240-207343/5), (LCS 240-207343/7), and (MB 240-207343/6), and found to be acceptable. As indicated in the reference method, sample analysis may proceed; however, any detection for the affected analytes is considered estimated.

The continuing calibration verification (CCV) analyzed in batch 240-207720 was outside the method criteria for the following analytes: trans-1,3-Dichloropropene and Dichloro-difluoromethane. An MRL standard at or below the reporting limit (RL) was analyzed with the

affected samples: F-VS-UST-P-2 (240-57930-2), (CCV 240-207720/5), (CCVIS 240-207720/4), (LCS 240-207720/6) and (MB 240-207720/8), and found to be acceptable. As indicated in the reference method, sample analysis may proceed; however, any detection for the affected analyte(s) is considered estimated.

The continuing calibration verification (CCV) analyzed in batch 240-207560 was outside the method criteria for the following analytes: 2-Hexanone, Acetone, 2-Methyl-2-propanol, 2-Chloroethyl Vinyl Ether, Chloromethane, 4-Methyl-2-pentanone, and 2-Butanone. An MRL standard at or below the reporting limit (RL) was analyzed with the affected samples: F-VS-UST-P-2 (240-57930-2), F-VS-UST-P-4 (240-57930-4), (CCV 240-207560/4), (CCVIS 240-207560/3), (LCS 240-207560/5), (LCS 240-207560/7) and (MB 240-207560/6) and found to be acceptable. As indicated in the reference method, sample analysis may proceed; however, any detection for the affected analytes is considered estimated.

An MS/MSD was analyzed in batch 240-207343. However due to poor purging or matrix interference the sample and MS/MSD could not be reported. Samples in the batch are: F-VS-UST-P-1 (240-57930-1), F-VS-UST-P-3 (240-57930-3), F-VB-UST-P-1 (240-57930-5), F-VB-UST-P-2 (240-57930-6), F-VB-UST-P-3 (240-57930-7), F-VB-UST-P-4 (240-57930-8), F-VB-UST-P-5 (240-57930-9), F-VB-UST-P-6 (240-57930-10), and F-VS-UST-P-DUP (240-57930-11).

Internal standard (ISTD) response for the following sample was outside control limits: F-VS-UST-P-4 (240-57930-4). The sample was re-extracted and/or re-analyzed with concurring results, and the original set of data has been reported.

Sample F-VS-UST-P-2 (240-57930-2) had a reanalysis that had different results than the original analysis. This is most likely due to non-homogeneity in the sample. Both runs are reported.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**VOLATILE ORGANIC COMPOUNDS (GCMS)**

Sample TB-111315 (240-57930-12) was analyzed for volatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8260C. The sample was analyzed on 11/17/2015.

Methylene Chloride was detected in method blank MB 240-207096/6 at a level exceeding the reporting limit. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

The laboratory control sample (LCS) for analytical batch 240-207096 recovered outside control limits for 1,2,3-Trichlorobenzene. This has been identified as a poor performing analyte when analyzed using this method; therefore, re-extraction/re-analysis was not performed: TB-111315 (240-57930-12).

2-Chloroethyl vinyl ether cannot be reliably recovered in an acid preserved sample: TB-111315 (240-57930-12).

The continuing calibration verification (CCV) for analytical batch 207096 exceeded control criteria for multiple compounds. The samples associated with this CCV were non-detects for the affected analytes. In accordance with the laboratory SOP, a low level CCV at the reporting limit (labeled as an MRL) was analyzed and the affected compounds were detected; therefore the data has been reported. No further corrective action was required: TB-111315 (240-57930-12).

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**SEMIVOLATILE ORGANIC COMPOUNDS (GCMS)**

Samples F-VS-UST-P-1 (240-57930-1), F-VS-UST-P-2 (240-57930-2), F-VS-UST-P-3 (240-57930-3), F-VS-UST-P-4 (240-57930-4), F-VB-UST-P-1 (240-57930-5), F-VB-UST-P-2 (240-57930-6), F-VB-UST-P-3 (240-57930-7), F-VB-UST-P-4 (240-57930-8), F-VB-UST-P-5 (240-57930-9), F-VB-UST-P-6 (240-57930-10) and F-VS-UST-P-DUP (240-57930-11) were analyzed for semivolatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8270D. The samples were prepared on 11/16/2015 and analyzed on 11/19/2015.

Bis(2-ethylhexyl) phthalate was detected in method blanks MB 240-206887/23-A and MB 240-206908/21-A at a level that was above the method detection limit but below the reporting limit. The value should be considered an estimate, and has been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

Sample F-VS-UST-P-3 (240-57930-3)[2.5X] required dilution prior to analysis. The reporting limits have been adjusted accordingly.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**GASOLINE RANGE ORGANICS**

Samples F-VS-UST-P-1 (240-57930-1), F-VS-UST-P-2 (240-57930-2), F-VS-UST-P-3 (240-57930-3), F-VS-UST-P-4 (240-57930-4), F-VB-UST-P-1 (240-57930-5), F-VB-UST-P-2 (240-57930-6), F-VB-UST-P-3 (240-57930-7), F-VB-UST-P-4 (240-57930-8), F-VB-UST-P-5 (240-57930-9), F-VB-UST-P-6 (240-57930-10) and F-VS-UST-P-DUP (240-57930-11) were analyzed for gasoline range organics in accordance with EPA SW-846 Method 8015C - GRO. The samples were prepared and analyzed on 11/16/2015, 11/17/2015 and 11/18/2015.

Gasoline Range Organics [C6 - C10] was detected in method blank MB 240-207078/1-A at a level exceeding the reporting limit. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**DIESEL RANGE ORGANICS (DRO)**
Samples F-VS-UST-P-1 (240-57930-1), F-VS-UST-P-2 (240-57930-2), F-VS-UST-P-3 (240-57930-3), F-VS-UST-P-4 (240-57930-4), F-VB-UST-P-1 (240-57930-5), F-VB-UST-P-2 (240-57930-6), F-VB-UST-P-3 (240-57930-7), F-VB-UST-P-4 (240-57930-8), F-VB-UST-P-5 (240-57930-9), F-VB-UST-P-6 (240-57930-10) and F-VS-UST-P-DUP (240-57930-11) were analyzed for diesel range organics (DRO) in accordance with EPA SW-846 Method 8015C - DRO. The samples were prepared on 11/16/2015 and 11/17/2015 and analyzed on 11/18/2015 and 11/19/2015.

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**POLYCHLORINATED BIPHENYLS (PCBS)**
Samples F-VS-UST-P-1 (240-57930-1), F-VS-UST-P-2 (240-57930-2), F-VS-UST-P-3 (240-57930-3), F-VS-UST-P-4 (240-57930-4), F-VB-UST-P-1 (240-57930-5), F-VB-UST-P-2 (240-57930-6), F-VB-UST-P-3 (240-57930-7), F-VB-UST-P-4 (240-57930-8), F-VB-UST-P-5 (240-57930-9), F-VB-UST-P-6 (240-57930-10) and F-VS-UST-P-DUP (240-57930-11) were analyzed for polychlorinated biphenyls (PCBs) in accordance with EPA SW-846 Method 8082A. The samples were prepared on 11/16/2015 and analyzed on 11/18/2015 and 11/19/2015.

Surrogates are added during the extraction process prior to dilution. When the sample dilution is 5X or greater, surrogate recoveries are diluted out and no corrective action is required. All of the samples in this data set analyzed for PCBs were subjected to the sulfuric acid cleanup procedure before instrumental analysis, per EPA Method 3665A.

The following samples required a tetrabutylammonium sulfite (TBA) clean-up to reduce matrix interferences caused by sulfur: F-VS-UST-P-1 (240-57930-1), F-VS-UST-P-2 (240-57930-2), F-VS-UST-P-3 (240-57930-3), F-VS-UST-P-4 (240-57930-4), F-VB-UST-P-2 (240-57930-6), F-VB-UST-P-3 (240-57930-7), F-VB-UST-P-6 (240-57930-10), F-VS-UST-P-DUP (240-57930-11), and (240-57880-E-1-A). Reagents: 2289919, 1848564, and 2283702.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**PERCENT SOLIDS**
Samples F-VS-UST-P-1 (240-57930-1), F-VS-UST-P-2 (240-57930-2), F-VS-UST-P-3 (240-57930-3), F-VS-UST-P-4 (240-57930-4), F-VB-UST-P-1 (240-57930-5), F-VB-UST-P-2 (240-57930-6), F-VB-UST-P-3 (240-57930-7), F-VB-UST-P-4 (240-57930-8), F-VB-UST-P-5 (240-57930-9), F-VB-UST-P-6 (240-57930-10) and F-VS-UST-P-DUP (240-57930-11) were analyzed for percent solids in accordance with EPA Method 160.3 MOD. The samples were analyzed on 11/16/2015.

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 206629

Lab Sample ID: STD5 240-206629/2 IC Client Sample ID:

Date Analyzed: 11/13/15 09:35 Lab File ID: 51113002.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.83 | Poor chromatography | gruberj | 11/13/15 09:55 |
| N-Nitrosodimethylamine | 3.27 | Poor chromatography | gruberj | 11/13/15 09:55 |
| Pyridine | 3.32 | Poor chromatography | gruberj | 11/13/15 09:55 |
| Benzoic acid | 6.93 | Poor chromatography | gruberj | 11/13/15 09:55 |

Lab Sample ID: STD4 240-206629/3 IC Client Sample ID:

Date Analyzed: 11/13/15 09:59 Lab File ID: 51113003.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.78 | Poor chromatography | gruberj | 11/13/15 11:07 |
| N-Nitrosodimethylamine | 3.27 | Poor chromatography | gruberj | 11/13/15 11:07 |
| Pyridine | 3.30 | Poor chromatography | gruberj | 11/13/15 11:07 |
| Benzoic acid | 6.91 | Poor chromatography | gruberj | 11/13/15 11:07 |

Lab Sample ID: STD3 240-206629/4 IC Client Sample ID:

Date Analyzed: 11/13/15 10:23 Lab File ID: 51113004.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.83 | Poor chromatography | gruberj | 11/13/15 11:09 |
| N-Nitrosodimethylamine | 3.27 | Poor chromatography | gruberj | 11/13/15 11:09 |
| Pyridine | 3.33 | Poor chromatography | gruberj | 11/13/15 11:09 |
| Benzoic acid | 6.90 | Poor chromatography | gruberj | 11/13/15 11:09 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 206629

Lab Sample ID: STD2 240-206629/5 IC Client Sample ID:

Date Analyzed: 11/13/15 10:48 Lab File ID: 51113005.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.78 | Poor chromatography | gruberj | 11/13/15 11:12 |
| N-Nitrosodimethylamine | 3.27 | Poor chromatography | gruberj | 11/13/15 11:12 |
| Pyridine | 3.33 | Poor chromatography | gruberj | 11/13/15 11:12 |

Lab Sample ID: STD1 240-206629/6 IC Client Sample ID:

Date Analyzed: 11/13/15 11:12 Lab File ID: 51113006.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Benzo[g,h,i]perylene | 16.94 | Poor chromatography | gruberj | 11/13/15 12:57 |

Lab Sample ID: STD8 240-206629/8 IC Client Sample ID:

Date Analyzed: 11/13/15 12:00 Lab File ID: 51113008.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.78 | Poor chromatography | gruberj | 11/13/15 13:02 |
| N-Nitrosodimethylamine | 3.27 | Poor chromatography | gruberj | 11/13/15 13:02 |
| Pyridine | 3.33 | Poor chromatography | gruberj | 11/13/15 13:02 |
| Caprolactam | 7.57 | Poor chromatography | gruberj | 11/13/15 13:02 |

Lab Sample ID: STD7 240-206629/9 IC Client Sample ID:

Date Analyzed: 11/13/15 12:24 Lab File ID: 51113009.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.78 | Poor chromatography | gruberj | 11/13/15 13:04 |
| N-Nitrosodimethylamine | 3.27 | Poor chromatography | gruberj | 11/13/15 13:04 |
| Pyridine | 3.30 | Poor chromatography | gruberj | 11/13/15 13:04 |
| Caprolactam | 7.56 | Poor chromatography | gruberj | 11/13/15 13:04 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 206629

Lab Sample ID: STD6 240-206629/10 ICIS Client Sample ID:

Date Analyzed: 11/13/15 12:48 Lab File ID: 51113010.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.78 | Poor chromatography | gruberj | 11/13/15 13:07 |
| N-Nitrosodimethylamine | 3.26 | Poor chromatography | gruberj | 11/13/15 13:07 |
| Pyridine | 3.30 | Poor chromatography | gruberj | 11/13/15 13:07 |
| Benzoic acid | 6.96 | Poor chromatography | gruberj | 11/13/15 13:07 |
| Caprolactam | 7.54 | Poor chromatography | gruberj | 11/13/15 13:07 |

Lab Sample ID: ICV 240-206629/11 Client Sample ID:

Date Analyzed: 11/13/15 13:12 Lab File ID: 51113011.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.78 | Poor chromatography | gruberj | 11/13/15 13:34 |

Lab Sample ID: STD9 240-206629/12 IC Client Sample ID:

Date Analyzed: 11/13/15 13:36 Lab File ID: 51113012.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.78 | Poor chromatography | gruberj | 11/13/15 13:53 |
| Pyridine | 3.30 | Poor chromatography | gruberj | 11/13/15 13:53 |
| Benzoic acid | 6.99 | Poor chromatography | gruberj | 11/13/15 13:53 |
| Caprolactam | 7.57 | Poor chromatography | gruberj | 11/13/15 13:53 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 207287

Lab Sample ID: CCV 240-207287/2 CCVIS Client Sample ID:

Date Analyzed: 11/18/15 08:26 Lab File ID: 51118002.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.75 | Poor chromatography | gruberj | 11/18/15 08:46 |
| N-Nitrosodimethylamine | 3.20 | Poor chromatography | gruberj | 11/18/15 08:46 |
| Pyridine | 3.27 | Poor chromatography | gruberj | 11/18/15 08:46 |
| Benzoic acid | 6.93 | Poor chromatography | gruberj | 11/18/15 08:46 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 207541

Lab Sample ID: CCV 240-207541/2 CCVIS Client Sample ID:

Date Analyzed: 11/19/15 09:50 Lab File ID: 51119002.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 2.72 | Poor chromatography | ulmanm | 11/19/15 10:12 |
| N-Nitrosodimethylamine | 3.17 | Poor chromatography | ulmanm | 11/19/15 10:12 |
| Pyridine | 3.23 | Poor chromatography | ulmanm | 11/19/15 10:12 |
| Benzoic acid | 6.93 | Poor chromatography | ulmanm | 11/19/15 10:12 |
| Caprolactam | 7.51 | Poor chromatography | ulmanm | 11/19/15 10:12 |

Lab Sample ID: 240-57930-2 Client Sample ID: F-VS-UST-P-2

Date Analyzed: 11/19/15 19:23 Lab File ID: 51119026.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Benzo[g,h,i]perylene | 16.85 | Poor chromatography | gruberj | 11/20/15 09:21 |

Lab Sample ID: 240-57930-11 Client Sample ID: F-VS-UST-P-DUP

Date Analyzed: 11/19/15 20:10 Lab File ID: 51119028.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Benzo[k]fluoranthene | 14.11 | Poor chromatography | gruberj | 11/20/15 09:26 |

Lab Sample ID: 240-57930-4 Client Sample ID: F-VS-UST-P-4

Date Analyzed: 11/19/15 20:34 Lab File ID: 51119029.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Benzo[k]fluoranthene | 14.12 | Poor chromatography | gruberj | 11/20/15 09:28 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A4HP10 Analysis Batch Number: 207541

Lab Sample ID: 240-57930-3 Client Sample ID: F-VS-UST-P-3

Date Analyzed: 11/19/15 20:57 Lab File ID: 51119030.D GC Column: DB-624 ID: 0.18(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Benzo[k]fluoranthene | 14.11 | Poor chromatography | gruberj | 11/20/15 09:30 |

GC VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: AFID Analysis Batch Number: 194508

Lab Sample ID: STD1 240-194508/7 IC Client Sample ID:

Date Analyzed: 08/24/15 13:59 Lab File ID: A0082107.D GC Column: RTX-VRX ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Trifluorotoluene (Surr) | 10.34 | Baseline Smoothing | bosworthh | 08/25/15 10:02 |
| Gasoline Range Organics [C6 - C10] | 11.46 | Baseline Smoothing | bosworthh | 08/25/15 10:02 |
| C6-C12 | 12.73 | Baseline Smoothing | bosworthh | 08/25/15 10:02 |

GC VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: AFID Analysis Batch Number: 207246

Lab Sample ID: 240-57930-3 Client Sample ID: F-VS-UST-P-3

Date Analyzed: 11/18/15 17:04 Lab File ID: A0111819.D GC Column: RTX-VRX ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Trifluorotoluene (Surr) | 10.37 | Baseline Smoothing | reedr | 11/19/15 06:16 |
| Gasoline Range Organics [C6 - C10] | 11.56 | Baseline Smoothing | reedr | 11/19/15 06:16 |

GC VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: YPID Analysis Batch Number: 206833

Lab Sample ID: MB 240-206878/1-A Client Sample ID:

Date Analyzed: 11/16/15 10:10 Lab File ID: Y0111609.D GC Column: RTX-VRX ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Gasoline Range Organics [C6 - C10] | 11.92 | Baseline Smoothing | reedr | 11/17/15 06:42 |

Lab Sample ID: 240-57930-1 Client Sample ID: F-VS-UST-P-1

Date Analyzed: 11/16/15 11:18 Lab File ID: Y0111611.D GC Column: RTX-VRX ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Trifluorotoluene (Surr) | 9.73 | Baseline Smoothing | reedr | 11/17/15 06:43 |
| Gasoline Range Organics [C6 - C10] | 11.92 | Baseline Smoothing | reedr | 11/17/15 06:43 |

DIESEL RANGE ORGANICS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP5R Analysis Batch Number: 207465

Lab Sample ID: CCV 240-207465/7 CCVRT Client Sample ID:

Date Analyzed: 11/18/15 21:17 Lab File ID: P5B11807.D GC Column: ID:

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| o-Terphenyl (Surr) | 9.83 | Other | bolgrind | 11/19/15 10:51 |

Lab Sample ID: CCV 240-207465/14 Client Sample ID:

Date Analyzed: 11/19/15 00:58 Lab File ID: P5B11814.D GC Column: ID:

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| o-Terphenyl (Surr) | 9.83 | Other | bolgrind | 11/19/15 10:56 |

Lab Sample ID: 240-57930-11 MS Client Sample ID: F-VS-UST-P-DUP MS

Date Analyzed: 11/19/15 04:38 Lab File ID: P5B11821.D GC Column: ID:

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| o-Terphenyl (Surr) | 9.83 | Other | bolgrind | 11/19/15 10:56 |

Lab Sample ID: 240-57930-11 MSD Client Sample ID: F-VS-UST-P-DUP MSD

Date Analyzed: 11/19/15 05:10 Lab File ID: P5B11822.D GC Column: ID:

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| o-Terphenyl (Surr) | 9.83 | Other | bolgrind | 11/19/15 10:56 |

Lab Sample ID: CCV 240-207465/23 Client Sample ID:

Date Analyzed: 11/19/15 05:42 Lab File ID: P5B11823.D GC Column: ID:

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| o-Terphenyl (Surr) | 9.83 | Other | bolgrind | 11/19/15 10:57 |

PCBS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP12 Analysis Batch Number: 207007

Lab Sample ID: STD05 240-207007/7 IC Client Sample ID:

Date Analyzed: 11/16/15 21:09 Lab File ID: P1200042.D GC Column: CLP-1 ID: 0.53(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| PCB-1262 Peak 1 | 7.95 | Baseline Smoothing | hassl | 11/17/15 06:39 |
| PCB-1262 Peak 2 | 8.75 | Baseline Smoothing | hassl | 11/17/15 06:39 |
| PCB-1262 Peak 3 | 9.03 | Baseline Smoothing | hassl | 11/17/15 06:39 |
| PCB-1262 Peak 4 | 9.40 | Baseline Smoothing | hassl | 11/17/15 06:39 |
| PCB-1262 Peak 5 | 9.77 | Baseline Smoothing | hassl | 11/17/15 06:39 |

Lab Sample ID: STD1 240-207007/8 IC Client Sample ID:

Date Analyzed: 11/16/15 21:27 Lab File ID: P1200043.D GC Column: CLP-1 ID: 0.53(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| PCB-1262 Peak 1 | 7.95 | Baseline Smoothing | hassl | 11/17/15 06:40 |
| PCB-1262 Peak 2 | 8.75 | Baseline Smoothing | hassl | 11/17/15 06:40 |
| PCB-1262 Peak 3 | 9.03 | Baseline Smoothing | hassl | 11/17/15 06:40 |
| PCB-1262 Peak 4 | 9.40 | Baseline Smoothing | hassl | 11/17/15 06:40 |
| PCB-1262 Peak 5 | 9.77 | Baseline Smoothing | hassl | 11/17/15 06:40 |

Lab Sample ID: STD02 240-207007/30 IC Client Sample ID:

Date Analyzed: 11/17/15 04:05 Lab File ID: P1200065.D GC Column: CLP-1 ID: 0.53(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| PCB-1016 Peak 1 | 4.61 | Baseline Smoothing | hassl | 11/17/15 06:42 |
| PCB-1016 Peak 2 | 5.11 | Baseline Smoothing | hassl | 11/17/15 06:42 |
| PCB-1016 Peak 3 | 5.74 | Baseline Smoothing | hassl | 11/17/15 06:42 |
| PCB-1016 Peak 4 | 5.91 | Baseline Smoothing | hassl | 11/17/15 06:42 |
| PCB-1016 Peak 5 | 6.45 | Baseline Smoothing | hassl | 11/17/15 06:42 |

PCBS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP12 Analysis Batch Number: 207007

Lab Sample ID: ICV 240-207007/39 Client Sample ID:

Date Analyzed: 11/17/15 06:48 Lab File ID: P1200074.D GC Column: CLP-1 ID: 0.53(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| PCB-1260 Peak 1 | 7.95 | Baseline Smoothing | hassl | 11/17/15 08:38 |
| PCB-1260 Peak 2 | 8.29 | Baseline Smoothing | hassl | 11/17/15 08:38 |
| PCB-1260 Peak 3 | 8.61 | Baseline Smoothing | hassl | 11/17/15 08:38 |
| PCB-1260 Peak 4 | 9.40 | Baseline Smoothing | hassl | 11/17/15 08:38 |
| PCB-1260 Peak 5 | 9.72 | Baseline Smoothing | hassl | 11/17/15 08:38 |

# SAMPLE SUMMARY

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Lab Sample ID | Client Sample ID | Client Matrix | Date/Time Sampled | Date/Time Received |
|---|---|---|---|---|
| 240-57930-1 | F-VS-UST-P-1 | Solid | 11/13/2015 1030 | 11/14/2015 0945 |
| 240-57930-2 | F-VS-UST-P-2 | Solid | 11/13/2015 1035 | 11/14/2015 0945 |
| 240-57930-3 | F-VS-UST-P-3 | Solid | 11/13/2015 1040 | 11/14/2015 0945 |
| 240-57930-4 | F-VS-UST-P-4 | Solid | 11/13/2015 1045 | 11/14/2015 0945 |
| 240-57930-5 | F-VB-UST-P-1 | Solid | 11/13/2015 1000 | 11/14/2015 0945 |
| 240-57930-6 | F-VB-UST-P-2 | Solid | 11/13/2015 1005 | 11/14/2015 0945 |
| 240-57930-7 | F-VB-UST-P-3 | Solid | 11/13/2015 1010 | 11/14/2015 0945 |
| 240-57930-8 | F-VB-UST-P-4 | Solid | 11/13/2015 1015 | 11/14/2015 0945 |
| 240-57930-9 | F-VB-UST-P-5 | Solid | 11/13/2015 1020 | 11/14/2015 0945 |
| 240-57930-10 | F-VB-UST-P-6 | Solid | 11/13/2015 1025 | 11/14/2015 0945 |
| 240-57930-11 | F-VS-UST-P-DUP | Solid | 11/13/2015 0000 | 11/14/2015 0945 |
| 240-57930-12 | TB-111315 | Water | 11/13/2015 0000 | 11/14/2015 0945 |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-57930-1** | **F-VS-UST-P-1** | | | | | |
| 1,2,3-Trimethylbenzene | | 0.87 | J | 5.7 | ug/Kg | 8260C |
| 2-Butanone (MEK) | | 18 | J | 23 | ug/Kg | 8260C |
| Acetone | | 86 | | 23 | ug/Kg | 8260C |
| Benzene | | 1.8 | J | 5.7 | ug/Kg | 8260C |
| Isopropylbenzene | | 14 | | 5.7 | ug/Kg | 8260C |
| m-Xylene & p-Xylene | | 1.9 | J | 11 | ug/Kg | 8260C |
| Naphthalene | | 1.1 | J | 5.7 | ug/Kg | 8260C |
| n-Butylbenzene | | 7.1 | | 5.7 | ug/Kg | 8260C |
| N-Propylbenzene | | 20 | | 5.7 | ug/Kg | 8260C |
| o-Xylene | | 0.75 | J | 5.7 | ug/Kg | 8260C |
| sec-Butylbenzene | | 8.6 | | 5.7 | ug/Kg | 8260C |
| Toluene | | 1.2 | J B | 5.7 | ug/Kg | 8260C |
| Xylenes, Total | | 2.7 | J | 11 | ug/Kg | 8260C |
| tert-Butylbenzene | | 2.3 | J | 5.7 | ug/Kg | 8260C |
| 2-Methylnaphthalene | | 7.1 | J | 8.0 | ug/Kg | 8270D |
| Anthracene | | 4.6 | J | 8.0 | ug/Kg | 8270D |
| Benzaldehyde | | 17 | J | 120 | ug/Kg | 8270D |
| Benzo[a]anthracene | | 12 | | 8.0 | ug/Kg | 8270D |
| Benzo[a]pyrene | | 12 | | 8.0 | ug/Kg | 8270D |
| Benzo[b]fluoranthene | | 17 | | 8.0 | ug/Kg | 8270D |
| Benzo[g,h,i]perylene | | 8.7 | | 8.0 | ug/Kg | 8270D |
| Benzo[k]fluoranthene | | 9.7 | | 8.0 | ug/Kg | 8270D |
| Bis(2-ethylhexyl) phthalate | | 23 | J B | 84 | ug/Kg | 8270D |
| Chrysene | | 18 | | 8.0 | ug/Kg | 8270D |
| Dibenzofuran | | 14 | J | 60 | ug/Kg | 8270D |
| Fluoranthene | | 27 | | 8.0 | ug/Kg | 8270D |
| Fluorene | | 4.8 | J | 8.0 | ug/Kg | 8270D |
| Indeno[1,2,3-cd]pyrene | | 7.3 | J | 8.0 | ug/Kg | 8270D |
| Naphthalene | | 7.4 | J | 8.0 | ug/Kg | 8270D |
| Phenanthrene | | 20 | | 8.0 | ug/Kg | 8270D |
| Pyrene | | 23 | | 8.0 | ug/Kg | 8270D |
| Gasoline Range Organics [C6 - C10] | | 3500 | | 1100 | ug/Kg | 8015C |
| Total Petroleum Hydrocarbons (C10-C32) | | 80 | | 19 | mg/Kg | 8015C |
| Percent Solids | | 85 | | 0.10 | % | Moisture |
| Percent Moisture | | 15 | | 0.10 | % | Moisture |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-57930-2** | **F-VS-UST-P-2** | | | | | |
| 1,2,3-Trimethylbenzene | | 7.9 | | 5.2 | ug/Kg | 8260C |
| 1,2,4-Trimethylbenzene | | 12 | | 5.2 | ug/Kg | 8260C |
| 2-Butanone (MEK) | | 6.1 | J | 21 | ug/Kg | 8260C |
| p-Isopropyltoluene | | 0.85 | J | 5.2 | ug/Kg | 8260C |
| Acetone | | 26 | B | 21 | ug/Kg | 8260C |
| Ethylbenzene | | 9.1 | | 5.2 | ug/Kg | 8260C |
| Isopropylbenzene | | 8.5 | | 5.2 | ug/Kg | 8260C |
| m-Xylene & p-Xylene | | 36 | | 10 | ug/Kg | 8260C |
| Naphthalene | | 10 | | 5.2 | ug/Kg | 8260C |
| n-Butylbenzene | | 4.6 | J | 5.2 | ug/Kg | 8260C |
| N-Propylbenzene | | 9.7 | | 5.2 | ug/Kg | 8260C |
| o-Xylene | | 2.7 | J | 5.2 | ug/Kg | 8260C |
| sec-Butylbenzene | | 3.6 | J | 5.2 | ug/Kg | 8260C |
| Toluene | | 1.2 | J B | 5.2 | ug/Kg | 8260C |
| Xylenes, Total | | 39 | | 10 | ug/Kg | 8260C |
| tert-Butylbenzene | | 0.87 | J | 5.2 | ug/Kg | 8260C |
| 1,1'-Biphenyl | | 7.2 | J | 60 | ug/Kg | 8270D |
| 2-Methylnaphthalene | | 85 | | 7.9 | ug/Kg | 8270D |
| Acenaphthene | | 10 | | 7.9 | ug/Kg | 8270D |
| Acenaphthylene | | 4.6 | J | 7.9 | ug/Kg | 8270D |
| Acetophenone | | 24 | J | 120 | ug/Kg | 8270D |
| Anthracene | | 7.4 | J | 7.9 | ug/Kg | 8270D |
| Benzaldehyde | | 23 | J | 120 | ug/Kg | 8270D |
| Benzo[a]anthracene | | 14 | | 7.9 | ug/Kg | 8270D |
| Benzo[a]pyrene | | 12 | | 7.9 | ug/Kg | 8270D |
| Benzo[b]fluoranthene | | 17 | | 7.9 | ug/Kg | 8270D |
| Benzo[g,h,i]perylene | | 10 | | 7.9 | ug/Kg | 8270D |
| Benzo[k]fluoranthene | | 8.9 | | 7.9 | ug/Kg | 8270D |
| Bis(2-ethylhexyl) phthalate | | 31 | J B | 83 | ug/Kg | 8270D |
| Chrysene | | 20 | | 7.9 | ug/Kg | 8270D |
| Dibenzofuran | | 11 | J | 60 | ug/Kg | 8270D |
| Fluoranthene | | 30 | | 7.9 | ug/Kg | 8270D |
| Fluorene | | 14 | | 7.9 | ug/Kg | 8270D |
| Indeno[1,2,3-cd]pyrene | | 9.0 | | 7.9 | ug/Kg | 8270D |
| Naphthalene | | 51 | | 7.9 | ug/Kg | 8270D |
| Phenanthrene | | 35 | | 7.9 | ug/Kg | 8270D |
| Pyrene | | 28 | | 7.9 | ug/Kg | 8270D |
| Gasoline Range Organics [C6 - C10] | | 5500 | B | 1200 | ug/Kg | 8015C |
| Total Petroleum Hydrocarbons (C10-C32) | | 130 | | 20 | mg/Kg | 8015C |
| Percent Solids | | 85 | | 0.10 | % | Moisture |
| Percent Moisture | | 15 | | 0.10 | % | Moisture |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-57930-3** | **F-VS-UST-P-3** | | | | | |
| 1,2,3-Trimethylbenzene | | 1.5 | J | 5.3 | ug/Kg | 8260C |
| 1,2,4-Trimethylbenzene | | 1.6 | J | 5.3 | ug/Kg | 8260C |
| 2-Butanone (MEK) | | 2.4 | J | 21 | ug/Kg | 8260C |
| Acetone | | 12 | J | 21 | ug/Kg | 8260C |
| Naphthalene | | 1.7 | J | 5.3 | ug/Kg | 8260C |
| Toluene | | 0.43 | J B | 5.3 | ug/Kg | 8260C |
| 1,1'-Biphenyl | | 85 | J | 150 | ug/Kg | 8270D |
| 2-Methylnaphthalene | | 360 | | 20 | ug/Kg | 8270D |
| Acenaphthene | | 350 | | 20 | ug/Kg | 8270D |
| Acenaphthylene | | 28 | | 20 | ug/Kg | 8270D |
| Anthracene | | 590 | | 20 | ug/Kg | 8270D |
| Benzo[a]anthracene | | 960 | | 20 | ug/Kg | 8270D |
| Benzo[a]pyrene | | 700 | | 20 | ug/Kg | 8270D |
| Benzo[b]fluoranthene | | 1100 | | 20 | ug/Kg | 8270D |
| Benzo[g,h,i]perylene | | 240 | | 20 | ug/Kg | 8270D |
| Benzo[k]fluoranthene | | 440 | | 20 | ug/Kg | 8270D |
| Carbazole | | 150 | | 150 | ug/Kg | 8270D |
| Chrysene | | 1000 | | 20 | ug/Kg | 8270D |
| Dibenz(a,h)anthracene | | 69 | | 20 | ug/Kg | 8270D |
| Dibenzofuran | | 330 | | 150 | ug/Kg | 8270D |
| Fluoranthene | | 2500 | | 20 | ug/Kg | 8270D |
| Fluorene | | 350 | | 20 | ug/Kg | 8270D |
| Indeno[1,2,3-cd]pyrene | | 250 | | 20 | ug/Kg | 8270D |
| Naphthalene | | 460 | | 20 | ug/Kg | 8270D |
| Phenanthrene | | 1900 | | 20 | ug/Kg | 8270D |
| Pyrene | | 1800 | | 20 | ug/Kg | 8270D |
| Gasoline Range Organics [C6 - C10] | | 110 | J | 120 | ug/Kg | 8015C |
| Total Petroleum Hydrocarbons (C10-C32) | | 42 | | 20 | mg/Kg | 8015C |
| Percent Solids | | 85 | | 0.10 | % | Moisture |
| Percent Moisture | | 15 | | 0.10 | % | Moisture |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-57930-4** | **F-VS-UST-P-4** | | | | | |
| 2-Butanone (MEK) | | 4.4 | J | 23 | ug/Kg | 8260C |
| Acetone | | 21 | J | 23 | ug/Kg | 8260C |
| Naphthalene | | 0.88 | J * | 5.7 | ug/Kg | 8260C |
| Toluene | | 0.47 | J B | 5.7 | ug/Kg | 8260C |
| 1,1'-Biphenyl | | 48 | J | 60 | ug/Kg | 8270D |
| 2-Methylnaphthalene | | 220 | | 7.9 | ug/Kg | 8270D |
| Acenaphthene | | 110 | | 7.9 | ug/Kg | 8270D |
| Acenaphthylene | | 23 | | 7.9 | ug/Kg | 8270D |
| Anthracene | | 160 | | 7.9 | ug/Kg | 8270D |
| Benzaldehyde | | 17 | J | 120 | ug/Kg | 8270D |
| Benzo[a]anthracene | | 200 | | 7.9 | ug/Kg | 8270D |
| Benzo[a]pyrene | | 100 | | 7.9 | ug/Kg | 8270D |
| Benzo[b]fluoranthene | | 240 | | 7.9 | ug/Kg | 8270D |
| Benzo[g,h,i]perylene | | 41 | | 7.9 | ug/Kg | 8270D |
| Benzo[k]fluoranthene | | 71 | | 7.9 | ug/Kg | 8270D |
| Bis(2-ethylhexyl) phthalate | | 33 | J B | 83 | ug/Kg | 8270D |
| Chrysene | | 260 | | 7.9 | ug/Kg | 8270D |
| Dibenz(a,h)anthracene | | 19 | | 7.9 | ug/Kg | 8270D |
| Dibenzofuran | | 170 | | 60 | ug/Kg | 8270D |
| Fluoranthene | | 560 | | 7.9 | ug/Kg | 8270D |
| Fluorene | | 120 | | 7.9 | ug/Kg | 8270D |
| Indeno[1,2,3-cd]pyrene | | 40 | | 7.9 | ug/Kg | 8270D |
| Naphthalene | | 290 | | 7.9 | ug/Kg | 8270D |
| Phenanthrene | | 550 | | 7.9 | ug/Kg | 8270D |
| Pyrene | | 420 | | 7.9 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 120 | | 20 | mg/Kg | 8015C |
| Percent Solids | | 83 | | 0.10 | % | Moisture |
| Percent Moisture | | 17 | | 0.10 | % | Moisture |
| **240-57930-5** | **F-VB-UST-P-1** | | | | | |
| 2-Butanone (MEK) | | 1.3 | J | 23 | ug/Kg | 8260C |
| Acetone | | 8.2 | J | 23 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 25 | J B | 82 | ug/Kg | 8270D |
| Percent Solids | | 85 | | 0.10 | % | Moisture |
| Percent Moisture | | 15 | | 0.10 | % | Moisture |
| **240-57930-6** | **F-VB-UST-P-2** | | | | | |
| 2-Butanone (MEK) | | 5.4 | J | 22 | ug/Kg | 8260C |
| Acetone | | 70 | | 22 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 38 | J B | 83 | ug/Kg | 8270D |
| Percent Solids | | 84 | | 0.10 | % | Moisture |
| Percent Moisture | | 16 | | 0.10 | % | Moisture |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-57930-7** | **F-VB-UST-P-3** | | | | | |
| 2-Butanone (MEK) | | 4.9 | J | 23 | ug/Kg | 8260C |
| Acetone | | 66 | | 23 | ug/Kg | 8260C |
| Benzaldehyde | | 45 | J | 120 | ug/Kg | 8270D |
| Bis(2-ethylhexyl) phthalate | | 55 | J B | 86 | ug/Kg | 8270D |
| Gasoline Range Organics [C6 - C10] | | 60 | J | 120 | ug/Kg | 8015C |
| Percent Solids | | 81 | | 0.10 | % | Moisture |
| Percent Moisture | | 19 | | 0.10 | % | Moisture |
| **240-57930-8** | **F-VB-UST-P-4** | | | | | |
| 2-Butanone (MEK) | | 1.9 | J | 21 | ug/Kg | 8260C |
| Acetone | | 16 | J | 21 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 30 | J B | 84 | ug/Kg | 8270D |
| Percent Solids | | 84 | | 0.10 | % | Moisture |
| Percent Moisture | | 16 | | 0.10 | % | Moisture |
| **240-57930-9** | **F-VB-UST-P-5** | | | | | |
| 2-Butanone (MEK) | | 2.0 | J | 22 | ug/Kg | 8260C |
| Acetone | | 11 | J | 22 | ug/Kg | 8260C |
| m-Xylene & p-Xylene | | 1.4 | J | 11 | ug/Kg | 8260C |
| Toluene | | 0.75 | J B | 5.4 | ug/Kg | 8260C |
| Xylenes, Total | | 1.4 | J | 11 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 62 | J B | 85 | ug/Kg | 8270D |
| Percent Solids | | 83 | | 0.10 | % | Moisture |
| Percent Moisture | | 17 | | 0.10 | % | Moisture |
| **240-57930-10** | **F-VB-UST-P-6** | | | | | |
| 2-Butanone (MEK) | | 4.1 | J | 23 | ug/Kg | 8260C |
| Acetone | | 57 | | 23 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 31 | J B | 82 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 8.0 | J | 20 | mg/Kg | 8015C |
| Percent Solids | | 85 | | 0.10 | % | Moisture |
| Percent Moisture | | 15 | | 0.10 | % | Moisture |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-57930-11** | **F-VS-UST-P-DUP** | | | | | |
| 2-Butanone (MEK) | | 6.4 | J | 21 | ug/Kg | 8260C |
| Acetone | | 34 | | 21 | ug/Kg | 8260C |
| Isopropylbenzene | | 2.8 | J | 5.2 | ug/Kg | 8260C |
| n-Butylbenzene | | 1.7 | J | 5.2 | ug/Kg | 8260C |
| N-Propylbenzene | | 4.4 | J | 5.2 | ug/Kg | 8260C |
| sec-Butylbenzene | | 2.0 | J | 5.2 | ug/Kg | 8260C |
| Toluene | | 0.31 | J B | 5.2 | ug/Kg | 8260C |
| 2-Methylnaphthalene | | 7.8 | | 7.8 | ug/Kg | 8270D |
| Acenaphthylene | | 4.0 | J | 7.8 | ug/Kg | 8270D |
| Acetophenone | | 17 | J | 120 | ug/Kg | 8270D |
| Anthracene | | 5.4 | J | 7.8 | ug/Kg | 8270D |
| Benzaldehyde | | 25 | J | 120 | ug/Kg | 8270D |
| Benzo[a]anthracene | | 16 | | 7.8 | ug/Kg | 8270D |
| Benzo[a]pyrene | | 17 | | 7.8 | ug/Kg | 8270D |
| Benzo[b]fluoranthene | | 30 | | 7.8 | ug/Kg | 8270D |
| Benzo[g,h,i]perylene | | 10 | | 7.8 | ug/Kg | 8270D |
| Benzo[k]fluoranthene | | 9.9 | | 7.8 | ug/Kg | 8270D |
| Chrysene | | 22 | | 7.8 | ug/Kg | 8270D |
| Dibenzofuran | | 16 | J | 58 | ug/Kg | 8270D |
| Fluoranthene | | 37 | | 7.8 | ug/Kg | 8270D |
| Fluorene | | 6.1 | J | 7.8 | ug/Kg | 8270D |
| Indeno[1,2,3-cd]pyrene | | 9.2 | | 7.8 | ug/Kg | 8270D |
| Naphthalene | | 12 | | 7.8 | ug/Kg | 8270D |
| Phenanthrene | | 27 | | 7.8 | ug/Kg | 8270D |
| Pyrene | | 30 | | 7.8 | ug/Kg | 8270D |
| Gasoline Range Organics [C6 - C10] | | 1800 | | 1200 | ug/Kg | 8015C |
| Total Petroleum Hydrocarbons (C10-C32) | | 51 | | 20 | mg/Kg | 8015C |
| Percent Solids | | 84 | | 0.10 | % | Moisture |
| Percent Moisture | | 16 | | 0.10 | % | Moisture |

# METHOD SUMMARY

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Description | Lab Location | Method | Preparation Method |
|---|---|---|---|
| **Matrix: Solid** | | | |
| Volatile Organic Compounds by GC/MS | TAL CAN | SW846 8260C | |
| Purge and Trap | TAL CAN | | SW846 5030C |
| Semivolatile Organic Compounds (GC/MS) | TAL CAN | SW846 8270D | |
| Soxhlet Extraction | TAL CAN | | SW846 3540C |
| Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics) | TAL CAN | SW846 8015C | |
| Purge and Trap | TAL CAN | | SW846 5030A |
| Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics) | TAL CAN | SW846 8015C | |
| Soxhlet Extraction | TAL CAN | | SW846 3540C |
| Polychlorinated Biphenyls (PCBs) by Gas Chromatography | TAL CAN | SW846 8082A | |
| Soxhlet Extraction | TAL CAN | | SW846 3540C |
| Percent Moisture | TAL CAN | EPA Moisture | |
| **Matrix: Water** | | | |
| Volatile Organic Compounds by GC/MS | TAL CAN | SW846 8260C | |
| Purge and Trap | TAL CAN | | SW846 5030C |

### Lab References:

TAL CAN = TestAmerica Canton

### Method References:

EPA = US Environmental Protection Agency

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

# METHOD / ANALYST SUMMARY

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Method | Analyst | Analyst ID |
|---|---|---|
| SW846 8260C | Macenczak, Steven | SAM |
| SW846 8260C | Quayle, Rick | RJQ |
| SW846 8270D | Gruber, John | JMG |
| SW846 8015C | Reed, Rachel T | RTR |
| SW846 8015C | Bolgrin, Deborah | DEB |
| SW846 8082A | Hass, Lori | LSH |
| EPA Moisture | Nguyen, Diem | DTN |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-1**

Lab Sample ID: 240-57930-1 Date Sampled: 11/13/2015 1030
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146551.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.18 g |
| Analysis Date: | 11/18/2015 1803 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.57 | U | 0.57 | 5.7 |
| 1,1,1-Trichloroethane | | 0.74 | U | 0.74 | 5.7 |
| 1,1,2,2-Tetrachloroethane | | 0.35 | U | 0.35 | 5.7 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.48 | U | 0.48 | 5.7 |
| 1,1-Dichloroethane | | 0.39 | U | 0.39 | 5.7 |
| 1,1-Dichloroethene | | 0.91 | U | 0.91 | 5.7 |
| 1,1-Dichloropropene | | 0.42 | U | 0.42 | 5.7 |
| 1,2,3-Trichlorobenzene | | 0.57 | U | 0.57 | 5.7 |
| 1,2,3-Trichloropropane | | 0.89 | U | 0.89 | 5.7 |
| 1,2,3-Trimethylbenzene | | 0.87 | J | 0.52 | 5.7 |
| 1,2,4-Trichlorobenzene | | 0.43 | U | 0.43 | 5.7 |
| 1,2,4-Trimethylbenzene | | 0.66 | U | 0.66 | 5.7 |
| 1,2-Dibromo-3-Chloropropane | | 3.1 | U | 3.1 | 11 |
| 1,2-Dichlorobenzene | | 0.39 | U | 0.39 | 5.7 |
| 1,2-Dichloroethane | | 0.52 | U | 0.52 | 5.7 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.7 |
| 1,3-Dichlorobenzene | | 0.78 | U | 0.78 | 5.7 |
| 1,3-Dichloropropane | | 0.45 | U | 0.45 | 5.7 |
| 1,4-Dichlorobenzene | | 0.32 | U | 0.32 | 5.7 |
| 2,2-Dichloropropane | | 0.50 | U | 0.50 | 5.7 |
| 2-Butanone (MEK) | | 18 | J | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.73 | U | 0.73 | 57 |
| 2-Chlorotoluene | | 0.34 | U | 0.34 | 5.7 |
| 2-Hexanone | | 0.77 | U | 0.77 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.7 |
| p-Isopropyltoluene | | 0.47 | U | 0.47 | 5.7 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 86 | | 6.5 | 23 |
| Benzene | | 1.8 | J | 1.0 | 5.7 |
| Bromobenzene | | 0.37 | U | 0.37 | 5.7 |
| Bromoform | | 0.27 | U | 0.27 | 5.7 |
| Bromomethane | | 0.48 | U | 0.48 | 5.7 |
| Carbon disulfide | | 0.67 | U | 0.67 | 5.7 |
| Carbon tetrachloride | | 0.74 | U | 0.74 | 5.7 |
| Chlorobenzene | | 0.60 | U | 0.60 | 5.7 |
| Dibromochloromethane | | 0.42 | U | 0.42 | 5.7 |
| Chloroethane | | 0.50 | U | 0.50 | 5.7 |
| Chloroform | | 0.41 | U | 0.41 | 5.7 |
| Chloromethane | | 0.90 | U | 0.90 | 5.7 |
| cis-1,2-Dichloroethene | | 0.32 | U | 0.32 | 5.7 |
| cis-1,3-Dichloropropene | | 0.81 | U | 0.81 | 5.7 |
| Dibromomethane | | 0.48 | U | 0.48 | 5.7 |
| Bromochloromethane | | 0.37 | U | 0.37 | 5.7 |
| Dichlorodifluoromethane | | 0.32 | U | 0.32 | 5.7 |
| Bromodichloromethane | | 0.33 | U | 0.33 | 5.7 |
| tert-Butyl alcohol (TBA) | | 6.5 | U | 6.5 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-1**

Lab Sample ID: 240-57930-1 Date Sampled: 11/13/2015 1030
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146551.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.18 g |
| Analysis Date: | 11/18/2015 1803 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.36 | U | 0.36 | 5.7 |
| Hexachlorobutadiene | | 0.45 | U | 0.45 | 5.7 |
| Diisopropyl ether | | 0.67 | U | 0.67 | 11 |
| Isopropylbenzene | | 14 | | 0.23 | 5.7 |
| Methyl tert-butyl ether | | 0.49 | U | 0.49 | 5.7 |
| Methylene Chloride | | 0.84 | U | 0.84 | 5.7 |
| m-Xylene & p-Xylene | | 1.9 | J | 0.48 | 11 |
| Naphthalene | | 1.1 | J | 0.40 | 5.7 |
| n-Butylbenzene | | 7.1 | | 0.67 | 5.7 |
| N-Propylbenzene | | 20 | | 0.68 | 5.7 |
| o-Xylene | | 0.75 | J | 0.65 | 5.7 |
| sec-Butylbenzene | | 8.6 | | 1.0 | 5.7 |
| Styrene | | 0.45 | U | 0.45 | 5.7 |
| Tert-amyl methyl ether | | 0.49 | U | 0.49 | 5.7 |
| Ethyl t-butyl ether | | 0.47 | U | 0.47 | 5.7 |
| Tetrachloroethene | | 0.91 | U | 0.91 | 5.7 |
| Toluene | | 1.2 | J B | 0.31 | 5.7 |
| trans-1,2-Dichloroethene | | 0.36 | U | 0.36 | 5.7 |
| trans-1,3-Dichloropropene | | 0.33 | U | 0.33 | 5.7 |
| Trichloroethene | | 0.43 | U | 0.43 | 5.7 |
| Trichlorofluoromethane | | 0.33 | U | 0.33 | 5.7 |
| Vinyl acetate | | 0.75 | U | 0.75 | 11 |
| Vinyl chloride | | 0.34 | U | 0.34 | 5.7 |
| Xylenes, Total | | 2.7 | J | 0.61 | 11 |
| tert-Butylbenzene | | 2.3 | J | 0.76 | 5.7 |
| 1,2-Dibromoethane | | 0.60 | U | 0.60 | 5.7 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 104 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 100 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 97 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 104 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207613 | Lab File ID: | 146583.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.74 g |
| Analysis Date: | 11/19/2015 1752 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.52 | U | 0.52 | 5.2 |
| 1,1,1-Trichloroethane | | 0.67 | U | 0.67 | 5.2 |
| 1,1,2,2-Tetrachloroethane | | 0.32 | U | 0.32 | 5.2 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.43 | U | 0.43 | 5.2 |
| 1,1-Dichloroethane | | 0.35 | U | 0.35 | 5.2 |
| 1,1-Dichloroethene | | 0.82 | U | 0.82 | 5.2 |
| 1,1-Dichloropropene | | 0.38 | U | 0.38 | 5.2 |
| 1,2,3-Trichlorobenzene | | 0.52 | U | 0.52 | 5.2 |
| 1,2,3-Trichloropropane | | 0.80 | U | 0.80 | 5.2 |
| 1,2,3-Trimethylbenzene | | 7.9 | | 0.47 | 5.2 |
| 1,2,4-Trichlorobenzene | | 0.39 | U | 0.39 | 5.2 |
| 1,2,4-Trimethylbenzene | | 12 | | 0.60 | 5.2 |
| 1,2-Dibromo-3-Chloropropane | | 2.8 | U | 2.8 | 10 |
| 1,2-Dichlorobenzene | | 0.35 | U | 0.35 | 5.2 |
| 1,2-Dichloroethane | | 0.47 | U | 0.47 | 5.2 |
| 1,2-Dichloropropane | | 0.14 | U | 0.14 | 5.2 |
| 1,3-Dichlorobenzene | | 0.71 | U | 0.71 | 5.2 |
| 1,3-Dichloropropane | | 0.41 | U | 0.41 | 5.2 |
| 1,4-Dichlorobenzene | | 0.29 | U | 0.29 | 5.2 |
| 2,2-Dichloropropane | | 0.45 | U | 0.45 | 5.2 |
| 2-Butanone (MEK) | | 6.1 | J | 1.2 | 21 |
| 2-Chloroethyl vinyl ether | | 0.66 | U | 0.66 | 52 |
| 2-Chlorotoluene | | 0.31 | U | 0.31 | 5.2 |
| 2-Hexanone | | 0.70 | U | 0.70 | 21 |
| 4-Chlorotoluene | | 0.27 | U | 0.27 | 5.2 |
| p-Isopropyltoluene | | 0.85 | J | 0.42 | 5.2 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 21 |
| Benzene | | 0.93 | U | 0.93 | 5.2 |
| Bromobenzene | | 0.34 | U | 0.34 | 5.2 |
| Bromoform | | 0.25 | U | 0.25 | 5.2 |
| Bromomethane | | 0.43 | U | 0.43 | 5.2 |
| Carbon disulfide | | 0.61 | U | 0.61 | 5.2 |
| Carbon tetrachloride | | 0.67 | U | 0.67 | 5.2 |
| Chlorobenzene | | 0.55 | U | 0.55 | 5.2 |
| Dibromochloromethane | | 0.38 | U | 0.38 | 5.2 |
| Chloroethane | | 0.45 | U | 0.45 | 5.2 |
| Chloroform | | 0.37 | U | 0.37 | 5.2 |
| Chloromethane | | 0.81 | U | 0.81 | 5.2 |
| cis-1,2-Dichloroethene | | 0.29 | U | 0.29 | 5.2 |
| cis-1,3-Dichloropropene | | 0.73 | U | 0.73 | 5.2 |
| Dibromomethane | | 0.43 | U | 0.43 | 5.2 |
| Bromochloromethane | | 0.34 | U | 0.34 | 5.2 |
| Dichlorodifluoromethane | | 0.29 | U | 0.29 | 5.2 |
| Bromodichloromethane | | 0.30 | U | 0.30 | 5.2 |
| tert-Butyl alcohol (TBA) | | 5.9 | U | 5.9 | 210 |
| Ethylbenzene | | 9.1 | | 0.33 | 5.2 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207613 | Lab File ID: | 146583.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.74 g |
| Analysis Date: | 11/19/2015 1752 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Hexachlorobutadiene | | 0.41 | U | 0.41 | 5.2 |
| Diisopropyl ether | | 0.61 | U | 0.61 | 10 |
| Isopropylbenzene | | 8.5 | | 0.21 | 5.2 |
| Methyl tert-butyl ether | | 0.44 | U | 0.44 | 5.2 |
| Methylene Chloride | | 0.76 | U | 0.76 | 5.2 |
| m-Xylene & p-Xylene | | 36 | | 0.43 | 10 |
| Naphthalene | | 10 | | 0.36 | 5.2 |
| n-Butylbenzene | | 4.6 | J | 0.61 | 5.2 |
| N-Propylbenzene | | 9.7 | | 0.62 | 5.2 |
| o-Xylene | | 2.7 | J | 0.59 | 5.2 |
| sec-Butylbenzene | | 3.6 | J | 0.94 | 5.2 |
| Styrene | | 0.41 | U | 0.41 | 5.2 |
| Tert-amyl methyl ether | | 0.44 | U | 0.44 | 5.2 |
| Ethyl t-butyl ether | | 0.42 | U | 0.42 | 5.2 |
| Tetrachloroethene | | 0.82 | U | 0.82 | 5.2 |
| Toluene | | 1.2 | J B | 0.28 | 5.2 |
| trans-1,2-Dichloroethene | | 0.33 | U | 0.33 | 5.2 |
| trans-1,3-Dichloropropene | | 0.30 | U | 0.30 | 5.2 |
| Trichloroethene | | 0.39 | U | 0.39 | 5.2 |
| Trichlorofluoromethane | | 0.30 | U | 0.30 | 5.2 |
| Vinyl acetate | | 0.68 | U | 0.68 | 10 |
| Vinyl chloride | | 0.31 | U | 0.31 | 5.2 |
| Xylenes, Total | | 39 | | 0.56 | 10 |
| tert-Butylbenzene | | 0.87 | J | 0.69 | 5.2 |
| 1,2-Dibromoethane | | 0.55 | U | 0.55 | 5.2 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 94 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 92 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 91 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 98 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207613 | Lab File ID: | 146584.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.09 g |
| Analysis Date: | 11/19/2015 1813 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.76 | U | 0.76 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.49 | U | 0.49 | 5.8 |
| 1,1-Dichloroethane | | 0.39 | U | 0.39 | 5.8 |
| 1,1-Dichloroethene | | 0.93 | U | 0.93 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.91 | U | 0.91 | 5.8 |
| 1,2,3-Trimethylbenzene | | 98 | | 0.53 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 170 | | 0.67 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.1 | U | 3.1 | 12 |
| 1,2-Dichlorobenzene | | 0.39 | U | 0.39 | 5.8 |
| 1,2-Dichloroethane | | 0.53 | U | 0.53 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.80 | U | 0.80 | 5.8 |
| 1,3-Dichloropropane | | 0.46 | U | 0.46 | 5.8 |
| 1,4-Dichlorobenzene | | 0.33 | U | 0.33 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 8.8 | J | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.74 | U | 0.74 | 58 |
| 2-Chlorotoluene | | 0.35 | U | 0.35 | 5.8 |
| 2-Hexanone | | 0.79 | U | 0.79 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.8 |
| p-Isopropyltoluene | | 6.5 | | 0.48 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Benzene | | 6.8 | | 1.0 | 5.8 |
| Bromobenzene | | 0.38 | U | 0.38 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.49 | U | 0.49 | 5.8 |
| Carbon disulfide | | 0.69 | U | 0.69 | 5.8 |
| Carbon tetrachloride | | 0.76 | U | 0.76 | 5.8 |
| Chlorobenzene | | 0.62 | U | 0.62 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.42 | U | 0.42 | 5.8 |
| Chloromethane | | 0.92 | U | 0.92 | 5.8 |
| cis-1,2-Dichloroethene | | 0.33 | U | 0.33 | 5.8 |
| cis-1,3-Dichloropropene | | 0.82 | U | 0.82 | 5.8 |
| Dibromomethane | | 0.49 | U | 0.49 | 5.8 |
| Bromochloromethane | | 0.38 | U | 0.38 | 5.8 |
| Dichlorodifluoromethane | | 0.33 | U | 0.33 | 5.8 |
| Bromodichloromethane | | 0.34 | U | 0.34 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.6 | U | 6.6 | 230 |
| Ethylbenzene | | 140 | | 0.37 | 5.8 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207613 | Lab File ID: | 146584.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.09 g |
| Analysis Date: | 11/19/2015 1813 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Hexachlorobutadiene | | 0.46 | U | 0.46 | 5.8 |
| Diisopropyl ether | | 0.69 | U | 0.69 | 12 |
| Isopropylbenzene | | 63 | | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.86 | U | 0.86 | 5.8 |
| m-Xylene & p-Xylene | | 520 | E | 0.49 | 12 |
| Naphthalene | | 17 | | 0.41 | 5.8 |
| n-Butylbenzene | | 11 | | 0.69 | 5.8 |
| N-Propylbenzene | | 55 | | 0.70 | 5.8 |
| o-Xylene | | 14 | | 0.66 | 5.8 |
| sec-Butylbenzene | | 12 | | 1.1 | 5.8 |
| Styrene | | 0.46 | U | 0.46 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.48 | U | 0.48 | 5.8 |
| Tetrachloroethene | | 0.93 | U | 0.93 | 5.8 |
| Toluene | | 4.0 | J B | 0.31 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.34 | U | 0.34 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.34 | U | 0.34 | 5.8 |
| Vinyl acetate | | 0.77 | U | 0.77 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 530 | | 0.63 | 12 |
| tert-Butylbenzene | | 4.4 | J | 0.78 | 5.8 |
| 1,2-Dibromoethane | | 0.62 | U | 0.62 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 103 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 92 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 105 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 95 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207720 | Instrument ID: | A3UX8 |
| Prep Method: | 5030C | Prep Batch: | 240-207709 | Lab File ID: | UX80893.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.61 g |
| Analysis Date: | 11/20/2015 1006 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/20/2015 0947 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Acetone | | 26 | B | 6.0 | 21 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 110 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 94 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 93 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 88 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-3**

Lab Sample ID: 240-57930-3 Date Sampled: 11/13/2015 1040
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146552.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/18/2015 1825 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.53 | U | 0.53 | 5.3 |
| 1,1,1-Trichloroethane | | 0.69 | U | 0.69 | 5.3 |
| 1,1,2,2-Tetrachloroethane | | 0.33 | U | 0.33 | 5.3 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.45 | U | 0.45 | 5.3 |
| 1,1-Dichloroethane | | 0.36 | U | 0.36 | 5.3 |
| 1,1-Dichloroethene | | 0.85 | U | 0.85 | 5.3 |
| 1,1-Dichloropropene | | 0.39 | U | 0.39 | 5.3 |
| 1,2,3-Trichlorobenzene | | 0.53 | U | 0.53 | 5.3 |
| 1,2,3-Trichloropropane | | 0.83 | U | 0.83 | 5.3 |
| 1,2,3-Trimethylbenzene | | 1.5 | J | 0.49 | 5.3 |
| 1,2,4-Trichlorobenzene | | 0.40 | U | 0.40 | 5.3 |
| 1,2,4-Trimethylbenzene | | 1.6 | J | 0.62 | 5.3 |
| 1,2-Dibromo-3-Chloropropane | | 2.9 | U | 2.9 | 11 |
| 1,2-Dichlorobenzene | | 0.36 | U | 0.36 | 5.3 |
| 1,2-Dichloroethane | | 0.49 | U | 0.49 | 5.3 |
| 1,2-Dichloropropane | | 0.15 | U | 0.15 | 5.3 |
| 1,3-Dichlorobenzene | | 0.73 | U | 0.73 | 5.3 |
| 1,3-Dichloropropane | | 0.43 | U | 0.43 | 5.3 |
| 1,4-Dichlorobenzene | | 0.30 | U | 0.30 | 5.3 |
| 2,2-Dichloropropane | | 0.47 | U | 0.47 | 5.3 |
| 2-Butanone (MEK) | | 2.4 | J | 1.2 | 21 |
| 2-Chloroethyl vinyl ether | | 0.68 | U | 0.68 | 53 |
| 2-Chlorotoluene | | 0.32 | U | 0.32 | 5.3 |
| 2-Hexanone | | 0.72 | U | 0.72 | 21 |
| 4-Chlorotoluene | | 0.28 | U | 0.28 | 5.3 |
| p-Isopropyltoluene | | 0.44 | U | 0.44 | 5.3 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 21 |
| Acetone | | 12 | J | 6.1 | 21 |
| Benzene | | 0.96 | U | 0.96 | 5.3 |
| Bromobenzene | | 0.35 | U | 0.35 | 5.3 |
| Bromoform | | 0.26 | U | 0.26 | 5.3 |
| Bromomethane | | 0.45 | U | 0.45 | 5.3 |
| Carbon disulfide | | 0.63 | U | 0.63 | 5.3 |
| Carbon tetrachloride | | 0.69 | U | 0.69 | 5.3 |
| Chlorobenzene | | 0.56 | U | 0.56 | 5.3 |
| Dibromochloromethane | | 0.39 | U | 0.39 | 5.3 |
| Chloroethane | | 0.47 | U | 0.47 | 5.3 |
| Chloroform | | 0.38 | U | 0.38 | 5.3 |
| Chloromethane | | 0.84 | U | 0.84 | 5.3 |
| cis-1,2-Dichloroethene | | 0.30 | U | 0.30 | 5.3 |
| cis-1,3-Dichloropropene | | 0.76 | U | 0.76 | 5.3 |
| Dibromomethane | | 0.45 | U | 0.45 | 5.3 |
| Bromochloromethane | | 0.35 | U | 0.35 | 5.3 |
| Dichlorodifluoromethane | | 0.30 | U | 0.30 | 5.3 |
| Bromodichloromethane | | 0.31 | U | 0.31 | 5.3 |
| tert-Butyl alcohol (TBA) | | 6.1 | U | 6.1 | 210 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-3**

Lab Sample ID: 240-57930-3 Date Sampled: 11/13/2015 1040
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146552.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/18/2015 1825 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.34 | U | 0.34 | 5.3 |
| Hexachlorobutadiene | | 0.43 | U | 0.43 | 5.3 |
| Diisopropyl ether | | 0.63 | U | 0.63 | 11 |
| Isopropylbenzene | | 0.21 | U | 0.21 | 5.3 |
| Methyl tert-butyl ether | | 0.46 | U | 0.46 | 5.3 |
| Methylene Chloride | | 0.79 | U | 0.79 | 5.3 |
| m-Xylene & p-Xylene | | 0.45 | U | 0.45 | 11 |
| Naphthalene | | 1.7 | J | 0.37 | 5.3 |
| n-Butylbenzene | | 0.63 | U | 0.63 | 5.3 |
| N-Propylbenzene | | 0.64 | U | 0.64 | 5.3 |
| o-Xylene | | 0.61 | U | 0.61 | 5.3 |
| sec-Butylbenzene | | 0.97 | U | 0.97 | 5.3 |
| Styrene | | 0.43 | U | 0.43 | 5.3 |
| Tert-amyl methyl ether | | 0.46 | U | 0.46 | 5.3 |
| Ethyl t-butyl ether | | 0.44 | U | 0.44 | 5.3 |
| Tetrachloroethene | | 0.85 | U | 0.85 | 5.3 |
| Toluene | | 0.43 | J B | 0.29 | 5.3 |
| trans-1,2-Dichloroethene | | 0.34 | U | 0.34 | 5.3 |
| trans-1,3-Dichloropropene | | 0.31 | U | 0.31 | 5.3 |
| Trichloroethene | | 0.40 | U | 0.40 | 5.3 |
| Trichlorofluoromethane | | 0.31 | U | 0.31 | 5.3 |
| Vinyl acetate | | 0.70 | U | 0.70 | 11 |
| Vinyl chloride | | 0.32 | U | 0.32 | 5.3 |
| Xylenes, Total | | 0.57 | U | 0.57 | 11 |
| tert-Butylbenzene | | 0.71 | U | 0.71 | 5.3 |
| 1,2-Dibromoethane | | 0.56 | U | 0.56 | 5.3 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 94 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 91 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 89 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 91 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-4**

Lab Sample ID: 240-57930-4 Date Sampled: 11/13/2015 1045
Client Matrix: Solid % Moisture: 16.9 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207613 | Lab File ID: | 146581.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.24 g |
| Analysis Date: | 11/19/2015 1708 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.57 | U | 0.57 | 5.7 |
| 1,1,1-Trichloroethane | | 0.75 | U | 0.75 | 5.7 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U * | 0.36 | 5.7 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.48 | U | 0.48 | 5.7 |
| 1,1-Dichloroethane | | 0.39 | U | 0.39 | 5.7 |
| 1,1-Dichloroethene | | 0.92 | U | 0.92 | 5.7 |
| 1,1-Dichloropropene | | 0.42 | U | 0.42 | 5.7 |
| 1,2,3-Trichlorobenzene | | 0.57 | U * | 0.57 | 5.7 |
| 1,2,3-Trichloropropane | | 0.90 | U * | 0.90 | 5.7 |
| 1,2,3-Trimethylbenzene | | 0.53 | U * | 0.53 | 5.7 |
| 1,2,4-Trichlorobenzene | | 0.44 | U * | 0.44 | 5.7 |
| 1,2,4-Trimethylbenzene | | 0.67 | U * | 0.67 | 5.7 |
| 1,2-Dibromo-3-Chloropropane | | 3.1 | U * | 3.1 | 11 |
| 1,2-Dichlorobenzene | | 0.39 | U * | 0.39 | 5.7 |
| 1,2-Dichloroethane | | 0.53 | U | 0.53 | 5.7 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.7 |
| 1,3-Dichlorobenzene | | 0.79 | U * | 0.79 | 5.7 |
| 1,3-Dichloropropane | | 0.46 | U | 0.46 | 5.7 |
| 1,4-Dichlorobenzene | | 0.32 | U * | 0.32 | 5.7 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.7 |
| 2-Butanone (MEK) | | 4.4 | J | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.73 | U | 0.73 | 57 |
| 2-Chlorotoluene | | 0.34 | U * | 0.34 | 5.7 |
| 2-Hexanone | | 0.78 | U | 0.78 | 23 |
| 4-Chlorotoluene | | 0.30 | U * | 0.30 | 5.7 |
| p-Isopropyltoluene | | 0.47 | U * | 0.47 | 5.7 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 21 | J | 6.5 | 23 |
| Benzene | | 1.0 | U | 1.0 | 5.7 |
| Bromobenzene | | 0.38 | U * | 0.38 | 5.7 |
| Bromoform | | 0.28 | U | 0.28 | 5.7 |
| Bromomethane | | 0.48 | U | 0.48 | 5.7 |
| Carbon disulfide | | 0.68 | U | 0.68 | 5.7 |
| Carbon tetrachloride | | 0.75 | U | 0.75 | 5.7 |
| Chlorobenzene | | 0.61 | U | 0.61 | 5.7 |
| Dibromochloromethane | | 0.42 | U | 0.42 | 5.7 |
| Chloroethane | | 0.51 | U | 0.51 | 5.7 |
| Chloroform | | 0.41 | U | 0.41 | 5.7 |
| Chloromethane | | 0.91 | U | 0.91 | 5.7 |
| cis-1,2-Dichloroethene | | 0.32 | U | 0.32 | 5.7 |
| cis-1,3-Dichloropropene | | 0.82 | U | 0.82 | 5.7 |
| Dibromomethane | | 0.48 | U | 0.48 | 5.7 |
| Bromochloromethane | | 0.38 | U | 0.38 | 5.7 |
| Dichlorodifluoromethane | | 0.32 | U | 0.32 | 5.7 |
| Bromodichloromethane | | 0.33 | U | 0.33 | 5.7 |
| tert-Butyl alcohol (TBA) | | 6.6 | U | 6.6 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-4**

Lab Sample ID: 240-57930-4 Date Sampled: 11/13/2015 1045
Client Matrix: Solid % Moisture: 16.9 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207613 | Lab File ID: | 146581.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.24 g |
| Analysis Date: | 11/19/2015 1708 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.37 | U | 0.37 | 5.7 |
| Hexachlorobutadiene | | 0.46 | U * | 0.46 | 5.7 |
| Diisopropyl ether | | 0.68 | U | 0.68 | 11 |
| Isopropylbenzene | | 0.23 | U | 0.23 | 5.7 |
| Methyl tert-butyl ether | | 0.49 | U | 0.49 | 5.7 |
| Methylene Chloride | | 0.85 | U | 0.85 | 5.7 |
| m-Xylene & p-Xylene | | 0.48 | U | 0.48 | 11 |
| Naphthalene | | 0.88 | J * | 0.40 | 5.7 |
| n-Butylbenzene | | 0.68 | U * | 0.68 | 5.7 |
| N-Propylbenzene | | 0.69 | U * | 0.69 | 5.7 |
| o-Xylene | | 0.65 | U | 0.65 | 5.7 |
| sec-Butylbenzene | | 1.0 | U * | 1.0 | 5.7 |
| Styrene | | 0.46 | U | 0.46 | 5.7 |
| Tert-amyl methyl ether | | 0.49 | U | 0.49 | 5.7 |
| Ethyl t-butyl ether | | 0.47 | U | 0.47 | 5.7 |
| Tetrachloroethene | | 0.92 | U | 0.92 | 5.7 |
| Toluene | | 0.47 | J B | 0.31 | 5.7 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.7 |
| trans-1,3-Dichloropropene | | 0.33 | U | 0.33 | 5.7 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.7 |
| Trichlorofluoromethane | | 0.33 | U | 0.33 | 5.7 |
| Vinyl acetate | | 0.76 | U | 0.76 | 11 |
| Vinyl chloride | | 0.34 | U | 0.34 | 5.7 |
| Xylenes, Total | | 0.62 | U | 0.62 | 11 |
| tert-Butylbenzene | | 0.77 | U * | 0.77 | 5.7 |
| 1,2-Dibromoethane | | 0.61 | U | 0.61 | 5.7 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 111 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 87 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 102 | * | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 85 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-1**

Lab Sample ID: 240-57930-5 Date Sampled: 11/13/2015 1000
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146556.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.06 g |
| Analysis Date: | 11/18/2015 1951 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.76 | U | 0.76 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.49 | U | 0.49 | 5.8 |
| 1,1-Dichloroethane | | 0.40 | U | 0.40 | 5.8 |
| 1,1-Dichloroethene | | 0.93 | U | 0.93 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.91 | U | 0.91 | 5.8 |
| 1,2,3-Trimethylbenzene | | 0.54 | U | 0.54 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 0.68 | U | 0.68 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.2 | U | 3.2 | 12 |
| 1,2-Dichlorobenzene | | 0.40 | U | 0.40 | 5.8 |
| 1,2-Dichloroethane | | 0.54 | U | 0.54 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.81 | U | 0.81 | 5.8 |
| 1,3-Dichloropropane | | 0.47 | U | 0.47 | 5.8 |
| 1,4-Dichlorobenzene | | 0.33 | U | 0.33 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 1.3 | J | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.75 | U | 0.75 | 58 |
| 2-Chlorotoluene | | 0.35 | U | 0.35 | 5.8 |
| 2-Hexanone | | 0.79 | U | 0.79 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.8 |
| p-Isopropyltoluene | | 0.48 | U | 0.48 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 8.2 | J | 6.7 | 23 |
| Benzene | | 1.1 | U | 1.1 | 5.8 |
| Bromobenzene | | 0.39 | U | 0.39 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.49 | U | 0.49 | 5.8 |
| Carbon disulfide | | 0.69 | U | 0.69 | 5.8 |
| Carbon tetrachloride | | 0.76 | U | 0.76 | 5.8 |
| Chlorobenzene | | 0.62 | U | 0.62 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.42 | U | 0.42 | 5.8 |
| Chloromethane | | 0.92 | U | 0.92 | 5.8 |
| cis-1,2-Dichloroethene | | 0.33 | U | 0.33 | 5.8 |
| cis-1,3-Dichloropropene | | 0.83 | U | 0.83 | 5.8 |
| Dibromomethane | | 0.49 | U | 0.49 | 5.8 |
| Bromochloromethane | | 0.39 | U | 0.39 | 5.8 |
| Dichlorodifluoromethane | | 0.33 | U | 0.33 | 5.8 |
| Bromodichloromethane | | 0.34 | U | 0.34 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.7 | U | 6.7 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-1**

Lab Sample ID: 240-57930-5 Date Sampled: 11/13/2015 1000
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146556.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.06 g |
| Analysis Date: | 11/18/2015 1951 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.37 | U | 0.37 | 5.8 |
| Hexachlorobutadiene | | 0.47 | U | 0.47 | 5.8 |
| Diisopropyl ether | | 0.69 | U | 0.69 | 12 |
| Isopropylbenzene | | 0.23 | U | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.86 | U | 0.86 | 5.8 |
| m-Xylene & p-Xylene | | 0.49 | U | 0.49 | 12 |
| Naphthalene | | 0.41 | U | 0.41 | 5.8 |
| n-Butylbenzene | | 0.69 | U | 0.69 | 5.8 |
| N-Propylbenzene | | 0.70 | U | 0.70 | 5.8 |
| o-Xylene | | 0.67 | U | 0.67 | 5.8 |
| sec-Butylbenzene | | 1.1 | U | 1.1 | 5.8 |
| Styrene | | 0.47 | U | 0.47 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.48 | U | 0.48 | 5.8 |
| Tetrachloroethene | | 0.93 | U | 0.93 | 5.8 |
| Toluene | | 0.32 | U | 0.32 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.34 | U | 0.34 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.34 | U | 0.34 | 5.8 |
| Vinyl acetate | | 0.77 | U | 0.77 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 0.63 | U | 0.63 | 12 |
| tert-Butylbenzene | | 0.78 | U | 0.78 | 5.8 |
| 1,2-Dibromoethane | | 0.62 | U | 0.62 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 89 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 89 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 78 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 90 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-2**

Lab Sample ID: 240-57930-6 Date Sampled: 11/13/2015 1005
Client Matrix: Solid % Moisture: 15.5 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146557.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.31 g |
| Analysis Date: | 11/18/2015 2012 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.56 | U | 0.56 | 5.6 |
| 1,1,1-Trichloroethane | | 0.72 | U | 0.72 | 5.6 |
| 1,1,2,2-Tetrachloroethane | | 0.35 | U | 0.35 | 5.6 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.47 | U | 0.47 | 5.6 |
| 1,1-Dichloroethane | | 0.38 | U | 0.38 | 5.6 |
| 1,1-Dichloroethene | | 0.89 | U | 0.89 | 5.6 |
| 1,1-Dichloropropene | | 0.41 | U | 0.41 | 5.6 |
| 1,2,3-Trichlorobenzene | | 0.56 | U | 0.56 | 5.6 |
| 1,2,3-Trichloropropane | | 0.87 | U | 0.87 | 5.6 |
| 1,2,3-Trimethylbenzene | | 0.51 | U | 0.51 | 5.6 |
| 1,2,4-Trichlorobenzene | | 0.42 | U | 0.42 | 5.6 |
| 1,2,4-Trimethylbenzene | | 0.65 | U | 0.65 | 5.6 |
| 1,2-Dibromo-3-Chloropropane | | 3.0 | U | 3.0 | 11 |
| 1,2-Dichlorobenzene | | 0.38 | U | 0.38 | 5.6 |
| 1,2-Dichloroethane | | 0.51 | U | 0.51 | 5.6 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.6 |
| 1,3-Dichlorobenzene | | 0.77 | U | 0.77 | 5.6 |
| 1,3-Dichloropropane | | 0.45 | U | 0.45 | 5.6 |
| 1,4-Dichlorobenzene | | 0.31 | U | 0.31 | 5.6 |
| 2,2-Dichloropropane | | 0.49 | U | 0.49 | 5.6 |
| 2-Butanone (MEK) | | 5.4 | J | 1.2 | 22 |
| 2-Chloroethyl vinyl ether | | 0.71 | U | 0.71 | 56 |
| 2-Chlorotoluene | | 0.33 | U | 0.33 | 5.6 |
| 2-Hexanone | | 0.76 | U | 0.76 | 22 |
| 4-Chlorotoluene | | 0.29 | U | 0.29 | 5.6 |
| p-Isopropyltoluene | | 0.46 | U | 0.46 | 5.6 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 22 |
| Acetone | | 70 | | 6.4 | 22 |
| Benzene | | 1.0 | U | 1.0 | 5.6 |
| Bromobenzene | | 0.37 | U | 0.37 | 5.6 |
| Bromoform | | 0.27 | U | 0.27 | 5.6 |
| Bromomethane | | 0.47 | U | 0.47 | 5.6 |
| Carbon disulfide | | 0.66 | U | 0.66 | 5.6 |
| Carbon tetrachloride | | 0.72 | U | 0.72 | 5.6 |
| Chlorobenzene | | 0.59 | U | 0.59 | 5.6 |
| Dibromochloromethane | | 0.41 | U | 0.41 | 5.6 |
| Chloroethane | | 0.49 | U | 0.49 | 5.6 |
| Chloroform | | 0.40 | U | 0.40 | 5.6 |
| Chloromethane | | 0.88 | U | 0.88 | 5.6 |
| cis-1,2-Dichloroethene | | 0.31 | U | 0.31 | 5.6 |
| cis-1,3-Dichloropropene | | 0.79 | U | 0.79 | 5.6 |
| Dibromomethane | | 0.47 | U | 0.47 | 5.6 |
| Bromochloromethane | | 0.37 | U | 0.37 | 5.6 |
| Dichlorodifluoromethane | | 0.31 | U | 0.31 | 5.6 |
| Bromodichloromethane | | 0.32 | U | 0.32 | 5.6 |
| tert-Butyl alcohol (TBA) | | 6.4 | U | 6.4 | 220 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-2**

Lab Sample ID: 240-57930-6 Date Sampled: 11/13/2015 1005
Client Matrix: Solid % Moisture: 15.5 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146557.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.31 g |
| Analysis Date: | 11/18/2015 2012 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.36 | U | 0.36 | 5.6 |
| Hexachlorobutadiene | | 0.45 | U | 0.45 | 5.6 |
| Diisopropyl ether | | 0.66 | U | 0.66 | 11 |
| Isopropylbenzene | | 0.22 | U | 0.22 | 5.6 |
| Methyl tert-butyl ether | | 0.48 | U | 0.48 | 5.6 |
| Methylene Chloride | | 0.82 | U | 0.82 | 5.6 |
| m-Xylene & p-Xylene | | 0.47 | U | 0.47 | 11 |
| Naphthalene | | 0.39 | U | 0.39 | 5.6 |
| n-Butylbenzene | | 0.66 | U | 0.66 | 5.6 |
| N-Propylbenzene | | 0.67 | U | 0.67 | 5.6 |
| o-Xylene | | 0.64 | U | 0.64 | 5.6 |
| sec-Butylbenzene | | 1.0 | U | 1.0 | 5.6 |
| Styrene | | 0.45 | U | 0.45 | 5.6 |
| Tert-amyl methyl ether | | 0.48 | U | 0.48 | 5.6 |
| Ethyl t-butyl ether | | 0.46 | U | 0.46 | 5.6 |
| Tetrachloroethene | | 0.89 | U | 0.89 | 5.6 |
| Toluene | | 0.30 | U | 0.30 | 5.6 |
| trans-1,2-Dichloroethene | | 0.36 | U | 0.36 | 5.6 |
| trans-1,3-Dichloropropene | | 0.32 | U | 0.32 | 5.6 |
| Trichloroethene | | 0.42 | U | 0.42 | 5.6 |
| Trichlorofluoromethane | | 0.32 | U | 0.32 | 5.6 |
| Vinyl acetate | | 0.74 | U | 0.74 | 11 |
| Vinyl chloride | | 0.33 | U | 0.33 | 5.6 |
| Xylenes, Total | | 0.60 | U | 0.60 | 11 |
| tert-Butylbenzene | | 0.75 | U | 0.75 | 5.6 |
| 1,2-Dibromoethane | | 0.59 | U | 0.59 | 5.6 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 89 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 95 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 79 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 100 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-3**

Lab Sample ID: 240-57930-7 Date Sampled: 11/13/2015 1010
Client Matrix: Solid % Moisture: 19.0 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146558.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.29 g |
| Analysis Date: | 11/18/2015 2033 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.76 | U | 0.76 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.49 | U | 0.49 | 5.8 |
| 1,1-Dichloroethane | | 0.40 | U | 0.40 | 5.8 |
| 1,1-Dichloroethene | | 0.93 | U | 0.93 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.91 | U | 0.91 | 5.8 |
| 1,2,3-Trimethylbenzene | | 0.54 | U | 0.54 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 0.68 | U | 0.68 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.2 | U | 3.2 | 12 |
| 1,2-Dichlorobenzene | | 0.40 | U | 0.40 | 5.8 |
| 1,2-Dichloroethane | | 0.54 | U | 0.54 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.81 | U | 0.81 | 5.8 |
| 1,3-Dichloropropane | | 0.47 | U | 0.47 | 5.8 |
| 1,4-Dichlorobenzene | | 0.33 | U | 0.33 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 4.9 | J | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.75 | U | 0.75 | 58 |
| 2-Chlorotoluene | | 0.35 | U | 0.35 | 5.8 |
| 2-Hexanone | | 0.79 | U | 0.79 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.8 |
| p-Isopropyltoluene | | 0.48 | U | 0.48 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 66 | | 6.7 | 23 |
| Benzene | | 1.1 | U | 1.1 | 5.8 |
| Bromobenzene | | 0.39 | U | 0.39 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.49 | U | 0.49 | 5.8 |
| Carbon disulfide | | 0.69 | U | 0.69 | 5.8 |
| Carbon tetrachloride | | 0.76 | U | 0.76 | 5.8 |
| Chlorobenzene | | 0.62 | U | 0.62 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.42 | U | 0.42 | 5.8 |
| Chloromethane | | 0.92 | U | 0.92 | 5.8 |
| cis-1,2-Dichloroethene | | 0.33 | U | 0.33 | 5.8 |
| cis-1,3-Dichloropropene | | 0.83 | U | 0.83 | 5.8 |
| Dibromomethane | | 0.49 | U | 0.49 | 5.8 |
| Bromochloromethane | | 0.39 | U | 0.39 | 5.8 |
| Dichlorodifluoromethane | | 0.33 | U | 0.33 | 5.8 |
| Bromodichloromethane | | 0.34 | U | 0.34 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.7 | U | 6.7 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-3**

Lab Sample ID: 240-57930-7 Date Sampled: 11/13/2015 1010
Client Matrix: Solid % Moisture: 19.0 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146558.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.29 g |
| Analysis Date: | 11/18/2015 2033 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.37 | U | 0.37 | 5.8 |
| Hexachlorobutadiene | | 0.47 | U | 0.47 | 5.8 |
| Diisopropyl ether | | 0.69 | U | 0.69 | 12 |
| Isopropylbenzene | | 0.23 | U | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.86 | U | 0.86 | 5.8 |
| m-Xylene & p-Xylene | | 0.49 | U | 0.49 | 12 |
| Naphthalene | | 0.41 | U | 0.41 | 5.8 |
| n-Butylbenzene | | 0.69 | U | 0.69 | 5.8 |
| N-Propylbenzene | | 0.70 | U | 0.70 | 5.8 |
| o-Xylene | | 0.67 | U | 0.67 | 5.8 |
| sec-Butylbenzene | | 1.1 | U | 1.1 | 5.8 |
| Styrene | | 0.47 | U | 0.47 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.48 | U | 0.48 | 5.8 |
| Tetrachloroethene | | 0.93 | U | 0.93 | 5.8 |
| Toluene | | 0.32 | U | 0.32 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.34 | U | 0.34 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.34 | U | 0.34 | 5.8 |
| Vinyl acetate | | 0.77 | U | 0.77 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 0.63 | U | 0.63 | 12 |
| tert-Butylbenzene | | 0.78 | U | 0.78 | 5.8 |
| 1,2-Dibromoethane | | 0.62 | U | 0.62 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 94 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 99 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 82 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 105 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-4**

Lab Sample ID: 240-57930-8 Date Sampled: 11/13/2015 1015
Client Matrix: Solid % Moisture: 16.1 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146559.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/18/2015 2055 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.54 | U | 0.54 | 5.4 |
| 1,1,1-Trichloroethane | | 0.70 | U | 0.70 | 5.4 |
| 1,1,2,2-Tetrachloroethane | | 0.33 | U | 0.33 | 5.4 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.45 | U | 0.45 | 5.4 |
| 1,1-Dichloroethane | | 0.37 | U | 0.37 | 5.4 |
| 1,1-Dichloroethene | | 0.86 | U | 0.86 | 5.4 |
| 1,1-Dichloropropene | | 0.40 | U | 0.40 | 5.4 |
| 1,2,3-Trichlorobenzene | | 0.54 | U | 0.54 | 5.4 |
| 1,2,3-Trichloropropane | | 0.84 | U | 0.84 | 5.4 |
| 1,2,3-Trimethylbenzene | | 0.49 | U | 0.49 | 5.4 |
| 1,2,4-Trichlorobenzene | | 0.41 | U | 0.41 | 5.4 |
| 1,2,4-Trimethylbenzene | | 0.62 | U | 0.62 | 5.4 |
| 1,2-Dibromo-3-Chloropropane | | 2.9 | U | 2.9 | 11 |
| 1,2-Dichlorobenzene | | 0.37 | U | 0.37 | 5.4 |
| 1,2-Dichloroethane | | 0.49 | U | 0.49 | 5.4 |
| 1,2-Dichloropropane | | 0.15 | U | 0.15 | 5.4 |
| 1,3-Dichlorobenzene | | 0.74 | U | 0.74 | 5.4 |
| 1,3-Dichloropropane | | 0.43 | U | 0.43 | 5.4 |
| 1,4-Dichlorobenzene | | 0.30 | U | 0.30 | 5.4 |
| 2,2-Dichloropropane | | 0.47 | U | 0.47 | 5.4 |
| 2-Butanone (MEK) | | 1.9 | J | 1.2 | 21 |
| 2-Chloroethyl vinyl ether | | 0.69 | U | 0.69 | 54 |
| 2-Chlorotoluene | | 0.32 | U | 0.32 | 5.4 |
| 2-Hexanone | | 0.73 | U | 0.73 | 21 |
| 4-Chlorotoluene | | 0.28 | U | 0.28 | 5.4 |
| p-Isopropyltoluene | | 0.44 | U | 0.44 | 5.4 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 21 |
| Acetone | | 16 | J | 6.1 | 21 |
| Benzene | | 0.97 | U | 0.97 | 5.4 |
| Bromobenzene | | 0.35 | U | 0.35 | 5.4 |
| Bromoform | | 0.26 | U | 0.26 | 5.4 |
| Bromomethane | | 0.45 | U | 0.45 | 5.4 |
| Carbon disulfide | | 0.63 | U | 0.63 | 5.4 |
| Carbon tetrachloride | | 0.70 | U | 0.70 | 5.4 |
| Chlorobenzene | | 0.57 | U | 0.57 | 5.4 |
| Dibromochloromethane | | 0.40 | U | 0.40 | 5.4 |
| Chloroethane | | 0.47 | U | 0.47 | 5.4 |
| Chloroform | | 0.39 | U | 0.39 | 5.4 |
| Chloromethane | | 0.85 | U | 0.85 | 5.4 |
| cis-1,2-Dichloroethene | | 0.30 | U | 0.30 | 5.4 |
| cis-1,3-Dichloropropene | | 0.76 | U | 0.76 | 5.4 |
| Dibromomethane | | 0.45 | U | 0.45 | 5.4 |
| Bromochloromethane | | 0.35 | U | 0.35 | 5.4 |
| Dichlorodifluoromethane | | 0.30 | U | 0.30 | 5.4 |
| Bromodichloromethane | | 0.31 | U | 0.31 | 5.4 |
| tert-Butyl alcohol (TBA) | | 6.1 | U | 6.1 | 210 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-4**

Lab Sample ID: 240-57930-8 Date Sampled: 11/13/2015 1015
Client Matrix: Solid % Moisture: 16.1 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146559.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/18/2015 2055 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.34 | U | 0.34 | 5.4 |
| Hexachlorobutadiene | | 0.43 | U | 0.43 | 5.4 |
| Diisopropyl ether | | 0.63 | U | 0.63 | 11 |
| Isopropylbenzene | | 0.21 | U | 0.21 | 5.4 |
| Methyl tert-butyl ether | | 0.46 | U | 0.46 | 5.4 |
| Methylene Chloride | | 0.79 | U | 0.79 | 5.4 |
| m-Xylene & p-Xylene | | 0.45 | U | 0.45 | 11 |
| Naphthalene | | 0.38 | U | 0.38 | 5.4 |
| n-Butylbenzene | | 0.63 | U | 0.63 | 5.4 |
| N-Propylbenzene | | 0.64 | U | 0.64 | 5.4 |
| o-Xylene | | 0.61 | U | 0.61 | 5.4 |
| sec-Butylbenzene | | 0.98 | U | 0.98 | 5.4 |
| Styrene | | 0.43 | U | 0.43 | 5.4 |
| Tert-amyl methyl ether | | 0.46 | U | 0.46 | 5.4 |
| Ethyl t-butyl ether | | 0.44 | U | 0.44 | 5.4 |
| Tetrachloroethene | | 0.86 | U | 0.86 | 5.4 |
| Toluene | | 0.29 | U | 0.29 | 5.4 |
| trans-1,2-Dichloroethene | | 0.34 | U | 0.34 | 5.4 |
| trans-1,3-Dichloropropene | | 0.31 | U | 0.31 | 5.4 |
| Trichloroethene | | 0.41 | U | 0.41 | 5.4 |
| Trichlorofluoromethane | | 0.31 | U | 0.31 | 5.4 |
| Vinyl acetate | | 0.71 | U | 0.71 | 11 |
| Vinyl chloride | | 0.32 | U | 0.32 | 5.4 |
| Xylenes, Total | | 0.58 | U | 0.58 | 11 |
| tert-Butylbenzene | | 0.72 | U | 0.72 | 5.4 |
| 1,2-Dibromoethane | | 0.57 | U | 0.57 | 5.4 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 96 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 103 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 89 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 113 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-5**

Lab Sample ID: 240-57930-9 Date Sampled: 11/13/2015 1020
Client Matrix: Solid % Moisture: 16.7 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146560.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/18/2015 2116 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.54 | U | 0.54 | 5.4 |
| 1,1,1-Trichloroethane | | 0.70 | U | 0.70 | 5.4 |
| 1,1,2,2-Tetrachloroethane | | 0.34 | U | 0.34 | 5.4 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.45 | U | 0.45 | 5.4 |
| 1,1-Dichloroethane | | 0.37 | U | 0.37 | 5.4 |
| 1,1-Dichloroethene | | 0.87 | U | 0.87 | 5.4 |
| 1,1-Dichloropropene | | 0.40 | U | 0.40 | 5.4 |
| 1,2,3-Trichlorobenzene | | 0.54 | U | 0.54 | 5.4 |
| 1,2,3-Trichloropropane | | 0.84 | U | 0.84 | 5.4 |
| 1,2,3-Trimethylbenzene | | 0.50 | U | 0.50 | 5.4 |
| 1,2,4-Trichlorobenzene | | 0.41 | U | 0.41 | 5.4 |
| 1,2,4-Trimethylbenzene | | 0.63 | U | 0.63 | 5.4 |
| 1,2-Dibromo-3-Chloropropane | | 2.9 | U | 2.9 | 11 |
| 1,2-Dichlorobenzene | | 0.37 | U | 0.37 | 5.4 |
| 1,2-Dichloroethane | | 0.50 | U | 0.50 | 5.4 |
| 1,2-Dichloropropane | | 0.15 | U | 0.15 | 5.4 |
| 1,3-Dichlorobenzene | | 0.75 | U | 0.75 | 5.4 |
| 1,3-Dichloropropane | | 0.43 | U | 0.43 | 5.4 |
| 1,4-Dichlorobenzene | | 0.30 | U | 0.30 | 5.4 |
| 2,2-Dichloropropane | | 0.48 | U | 0.48 | 5.4 |
| 2-Butanone (MEK) | | 2.0 | J | 1.2 | 22 |
| 2-Chloroethyl vinyl ether | | 0.69 | U | 0.69 | 54 |
| 2-Chlorotoluene | | 0.32 | U | 0.32 | 5.4 |
| 2-Hexanone | | 0.74 | U | 0.74 | 22 |
| 4-Chlorotoluene | | 0.28 | U | 0.28 | 5.4 |
| p-Isopropyltoluene | | 0.44 | U | 0.44 | 5.4 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 22 |
| Acetone | | 11 | J | 6.2 | 22 |
| Benzene | | 0.97 | U | 0.97 | 5.4 |
| Bromobenzene | | 0.36 | U | 0.36 | 5.4 |
| Bromoform | | 0.26 | U | 0.26 | 5.4 |
| Bromomethane | | 0.45 | U | 0.45 | 5.4 |
| Carbon disulfide | | 0.64 | U | 0.64 | 5.4 |
| Carbon tetrachloride | | 0.70 | U | 0.70 | 5.4 |
| Chlorobenzene | | 0.57 | U | 0.57 | 5.4 |
| Dibromochloromethane | | 0.40 | U | 0.40 | 5.4 |
| Chloroethane | | 0.48 | U | 0.48 | 5.4 |
| Chloroform | | 0.39 | U | 0.39 | 5.4 |
| Chloromethane | | 0.85 | U | 0.85 | 5.4 |
| cis-1,2-Dichloroethene | | 0.30 | U | 0.30 | 5.4 |
| cis-1,3-Dichloropropene | | 0.77 | U | 0.77 | 5.4 |
| Dibromomethane | | 0.45 | U | 0.45 | 5.4 |
| Bromochloromethane | | 0.36 | U | 0.36 | 5.4 |
| Dichlorodifluoromethane | | 0.30 | U | 0.30 | 5.4 |
| Bromodichloromethane | | 0.31 | U | 0.31 | 5.4 |
| tert-Butyl alcohol (TBA) | | 6.2 | U | 6.2 | 220 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-5**

Lab Sample ID: 240-57930-9 Date Sampled: 11/13/2015 1020
Client Matrix: Solid % Moisture: 16.7 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146560.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.55 g |
| Analysis Date: | 11/18/2015 2116 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.35 | U | 0.35 | 5.4 |
| Hexachlorobutadiene | | 0.43 | U | 0.43 | 5.4 |
| Diisopropyl ether | | 0.64 | U | 0.64 | 11 |
| Isopropylbenzene | | 0.22 | U | 0.22 | 5.4 |
| Methyl tert-butyl ether | | 0.47 | U | 0.47 | 5.4 |
| Methylene Chloride | | 0.80 | U | 0.80 | 5.4 |
| m-Xylene & p-Xylene | | 1.4 | J | 0.45 | 11 |
| Naphthalene | | 0.38 | U | 0.38 | 5.4 |
| n-Butylbenzene | | 0.64 | U | 0.64 | 5.4 |
| N-Propylbenzene | | 0.65 | U | 0.65 | 5.4 |
| o-Xylene | | 0.62 | U | 0.62 | 5.4 |
| sec-Butylbenzene | | 0.98 | U | 0.98 | 5.4 |
| Styrene | | 0.43 | U | 0.43 | 5.4 |
| Tert-amyl methyl ether | | 0.47 | U | 0.47 | 5.4 |
| Ethyl t-butyl ether | | 0.44 | U | 0.44 | 5.4 |
| Tetrachloroethene | | 0.87 | U | 0.87 | 5.4 |
| Toluene | | 0.75 | J B | 0.29 | 5.4 |
| trans-1,2-Dichloroethene | | 0.35 | U | 0.35 | 5.4 |
| trans-1,3-Dichloropropene | | 0.31 | U | 0.31 | 5.4 |
| Trichloroethene | | 0.41 | U | 0.41 | 5.4 |
| Trichlorofluoromethane | | 0.31 | U | 0.31 | 5.4 |
| Vinyl acetate | | 0.71 | U | 0.71 | 11 |
| Vinyl chloride | | 0.32 | U | 0.32 | 5.4 |
| Xylenes, Total | | 1.4 | J | 0.58 | 11 |
| tert-Butylbenzene | | 0.72 | U | 0.72 | 5.4 |
| 1,2-Dibromoethane | | 0.57 | U | 0.57 | 5.4 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 84 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 86 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 76 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 89 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-6**

Lab Sample ID: 240-57930-10 Date Sampled: 11/13/2015 1025
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146561.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.10 g |
| Analysis Date: | 11/18/2015 2138 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.75 | U | 0.75 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.49 | U | 0.49 | 5.8 |
| 1,1-Dichloroethane | | 0.39 | U | 0.39 | 5.8 |
| 1,1-Dichloroethene | | 0.93 | U | 0.93 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.90 | U | 0.90 | 5.8 |
| 1,2,3-Trimethylbenzene | | 0.53 | U | 0.53 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 0.67 | U | 0.67 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.1 | U | 3.1 | 12 |
| 1,2-Dichlorobenzene | | 0.39 | U | 0.39 | 5.8 |
| 1,2-Dichloroethane | | 0.53 | U | 0.53 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.80 | U | 0.80 | 5.8 |
| 1,3-Dichloropropane | | 0.46 | U | 0.46 | 5.8 |
| 1,4-Dichlorobenzene | | 0.32 | U | 0.32 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 4.1 | J | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.74 | U | 0.74 | 58 |
| 2-Chlorotoluene | | 0.35 | U | 0.35 | 5.8 |
| 2-Hexanone | | 0.79 | U | 0.79 | 23 |
| 4-Chlorotoluene | | 0.30 | U | 0.30 | 5.8 |
| p-Isopropyltoluene | | 0.47 | U | 0.47 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 57 | | 6.6 | 23 |
| Benzene | | 1.0 | U | 1.0 | 5.8 |
| Bromobenzene | | 0.38 | U | 0.38 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.49 | U | 0.49 | 5.8 |
| Carbon disulfide | | 0.68 | U | 0.68 | 5.8 |
| Carbon tetrachloride | | 0.75 | U | 0.75 | 5.8 |
| Chlorobenzene | | 0.61 | U | 0.61 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.42 | U | 0.42 | 5.8 |
| Chloromethane | | 0.91 | U | 0.91 | 5.8 |
| cis-1,2-Dichloroethene | | 0.32 | U | 0.32 | 5.8 |
| cis-1,3-Dichloropropene | | 0.82 | U | 0.82 | 5.8 |
| Dibromomethane | | 0.49 | U | 0.49 | 5.8 |
| Bromochloromethane | | 0.38 | U | 0.38 | 5.8 |
| Dichlorodifluoromethane | | 0.32 | U | 0.32 | 5.8 |
| Bromodichloromethane | | 0.34 | U | 0.34 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.6 | U | 6.6 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-6**

Lab Sample ID: 240-57930-10 Date Sampled: 11/13/2015 1025
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146561.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.10 g |
| Analysis Date: | 11/18/2015 2138 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1227 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.37 | U | 0.37 | 5.8 |
| Hexachlorobutadiene | | 0.46 | U | 0.46 | 5.8 |
| Diisopropyl ether | | 0.68 | U | 0.68 | 12 |
| Isopropylbenzene | | 0.23 | U | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.86 | U | 0.86 | 5.8 |
| m-Xylene & p-Xylene | | 0.49 | U | 0.49 | 12 |
| Naphthalene | | 0.40 | U | 0.40 | 5.8 |
| n-Butylbenzene | | 0.68 | U | 0.68 | 5.8 |
| N-Propylbenzene | | 0.69 | U | 0.69 | 5.8 |
| o-Xylene | | 0.66 | U | 0.66 | 5.8 |
| sec-Butylbenzene | | 1.1 | U | 1.1 | 5.8 |
| Styrene | | 0.46 | U | 0.46 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.47 | U | 0.47 | 5.8 |
| Tetrachloroethene | | 0.93 | U | 0.93 | 5.8 |
| Toluene | | 0.31 | U | 0.31 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.34 | U | 0.34 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.34 | U | 0.34 | 5.8 |
| Vinyl acetate | | 0.76 | U | 0.76 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 0.62 | U | 0.62 | 12 |
| tert-Butylbenzene | | 0.77 | U | 0.77 | 5.8 |
| 1,2-Dibromoethane | | 0.61 | U | 0.61 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 93 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 97 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 86 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 105 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-DUP**

Lab Sample ID: 240-57930-11 Date Sampled: 11/13/2015 0000
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146562.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.73 g |
| Analysis Date: | 11/18/2015 2159 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1230 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.52 | U | 0.52 | 5.2 |
| 1,1,1-Trichloroethane | | 0.67 | U | 0.67 | 5.2 |
| 1,1,2,2-Tetrachloroethane | | 0.32 | U | 0.32 | 5.2 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.43 | U | 0.43 | 5.2 |
| 1,1-Dichloroethane | | 0.35 | U | 0.35 | 5.2 |
| 1,1-Dichloroethene | | 0.83 | U | 0.83 | 5.2 |
| 1,1-Dichloropropene | | 0.38 | U | 0.38 | 5.2 |
| 1,2,3-Trichlorobenzene | | 0.52 | U | 0.52 | 5.2 |
| 1,2,3-Trichloropropane | | 0.81 | U | 0.81 | 5.2 |
| 1,2,3-Trimethylbenzene | | 0.48 | U | 0.48 | 5.2 |
| 1,2,4-Trichlorobenzene | | 0.39 | U | 0.39 | 5.2 |
| 1,2,4-Trimethylbenzene | | 0.60 | U | 0.60 | 5.2 |
| 1,2-Dibromo-3-Chloropropane | | 2.8 | U | 2.8 | 10 |
| 1,2-Dichlorobenzene | | 0.35 | U | 0.35 | 5.2 |
| 1,2-Dichloroethane | | 0.48 | U | 0.48 | 5.2 |
| 1,2-Dichloropropane | | 0.14 | U | 0.14 | 5.2 |
| 1,3-Dichlorobenzene | | 0.71 | U | 0.71 | 5.2 |
| 1,3-Dichloropropane | | 0.41 | U | 0.41 | 5.2 |
| 1,4-Dichlorobenzene | | 0.29 | U | 0.29 | 5.2 |
| 2,2-Dichloropropane | | 0.45 | U | 0.45 | 5.2 |
| 2-Butanone (MEK) | | 6.4 | J | 1.2 | 21 |
| 2-Chloroethyl vinyl ether | | 0.66 | U | 0.66 | 52 |
| 2-Chlorotoluene | | 0.31 | U | 0.31 | 5.2 |
| 2-Hexanone | | 0.70 | U | 0.70 | 21 |
| 4-Chlorotoluene | | 0.27 | U | 0.27 | 5.2 |
| p-Isopropyltoluene | | 0.42 | U | 0.42 | 5.2 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 21 |
| Acetone | | 34 | | 5.9 | 21 |
| Benzene | | 0.93 | U | 0.93 | 5.2 |
| Bromobenzene | | 0.34 | U | 0.34 | 5.2 |
| Bromoform | | 0.25 | U | 0.25 | 5.2 |
| Bromomethane | | 0.43 | U | 0.43 | 5.2 |
| Carbon disulfide | | 0.61 | U | 0.61 | 5.2 |
| Carbon tetrachloride | | 0.67 | U | 0.67 | 5.2 |
| Chlorobenzene | | 0.55 | U | 0.55 | 5.2 |
| Dibromochloromethane | | 0.38 | U | 0.38 | 5.2 |
| Chloroethane | | 0.45 | U | 0.45 | 5.2 |
| Chloroform | | 0.37 | U | 0.37 | 5.2 |
| Chloromethane | | 0.82 | U | 0.82 | 5.2 |
| cis-1,2-Dichloroethene | | 0.29 | U | 0.29 | 5.2 |
| cis-1,3-Dichloropropene | | 0.73 | U | 0.73 | 5.2 |
| Dibromomethane | | 0.43 | U | 0.43 | 5.2 |
| Bromochloromethane | | 0.34 | U | 0.34 | 5.2 |
| Dichlorodifluoromethane | | 0.29 | U | 0.29 | 5.2 |
| Bromodichloromethane | | 0.30 | U | 0.30 | 5.2 |
| tert-Butyl alcohol (TBA) | | 5.9 | U | 5.9 | 210 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-DUP**

Lab Sample ID: 240-57930-11 Date Sampled: 11/13/2015 0000
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-207393 | Lab File ID: | 146562.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.73 g |
| Analysis Date: | 11/18/2015 2159 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 1230 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.33 | U | 0.33 | 5.2 |
| Hexachlorobutadiene | | 0.41 | U | 0.41 | 5.2 |
| Diisopropyl ether | | 0.61 | U | 0.61 | 10 |
| Isopropylbenzene | | 2.8 | J | 0.21 | 5.2 |
| Methyl tert-butyl ether | | 0.44 | U | 0.44 | 5.2 |
| Methylene Chloride | | 0.76 | U | 0.76 | 5.2 |
| m-Xylene & p-Xylene | | 0.43 | U | 0.43 | 10 |
| Naphthalene | | 0.36 | U | 0.36 | 5.2 |
| n-Butylbenzene | | 1.7 | J | 0.61 | 5.2 |
| N-Propylbenzene | | 4.4 | J | 0.62 | 5.2 |
| o-Xylene | | 0.59 | U | 0.59 | 5.2 |
| sec-Butylbenzene | | 2.0 | J | 0.94 | 5.2 |
| Styrene | | 0.41 | U | 0.41 | 5.2 |
| Tert-amyl methyl ether | | 0.44 | U | 0.44 | 5.2 |
| Ethyl t-butyl ether | | 0.42 | U | 0.42 | 5.2 |
| Tetrachloroethene | | 0.83 | U | 0.83 | 5.2 |
| Toluene | | 0.31 | J B | 0.28 | 5.2 |
| trans-1,2-Dichloroethene | | 0.33 | U | 0.33 | 5.2 |
| trans-1,3-Dichloropropene | | 0.30 | U | 0.30 | 5.2 |
| Trichloroethene | | 0.39 | U | 0.39 | 5.2 |
| Trichlorofluoromethane | | 0.30 | U | 0.30 | 5.2 |
| Vinyl acetate | | 0.68 | U | 0.68 | 10 |
| Vinyl chloride | | 0.31 | U | 0.31 | 5.2 |
| Xylenes, Total | | 0.56 | U | 0.56 | 10 |
| tert-Butylbenzene | | 0.69 | U | 0.69 | 5.2 |
| 1,2-Dibromoethane | | 0.55 | U | 0.55 | 5.2 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 90 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 89 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 83 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 94 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: TB-111315**

Lab Sample ID: 240-57930-12 Date Sampled: 11/13/2015 0000
Client Matrix: Water Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Prep Method: | 5030C | Prep Batch: | N/A | Lab File ID: | UXX3077.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1451 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1451 | | | | |

| Analyte | Result (ug/L) | Qualifier | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.28 | U | 0.28 | 1.0 |
| 1,1,1-Trichloroethane | 0.44 | U | 0.44 | 1.0 |
| 1,1,2,2-Tetrachloroethane | 0.22 | U | 0.22 | 1.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloroethane | 0.30 | U | 0.30 | 1.0 |
| 1,1-Dichloroethene | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloropropene | 0.42 | U | 0.42 | 1.0 |
| 1,2,3-Trichlorobenzene | 0.37 | U * | 0.37 | 1.0 |
| 1,2,3-Trichloropropane | 0.44 | U | 0.44 | 1.0 |
| 1,2,3-Trimethylbenzene | 0.47 | U | 0.47 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.32 | U | 0.32 | 1.0 |
| 1,2,4-Trimethylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromo-3-Chloropropane | 0.82 | U | 0.82 | 2.0 |
| 1,2-Dichlorobenzene | 0.25 | U | 0.25 | 1.0 |
| 1,2-Dichloroethane | 0.23 | U | 0.23 | 1.0 |
| 1,2-Dichloropropane | 0.25 | U | 0.25 | 1.0 |
| 1,3-Dichlorobenzene | 0.19 | U | 0.19 | 1.0 |
| 1,3-Dichloropropane | 0.19 | U | 0.19 | 1.0 |
| 1,4-Dichlorobenzene | 0.27 | U | 0.27 | 1.0 |
| 2,2-Dichloropropane | 0.26 | U | 0.26 | 1.0 |
| 2-Butanone (MEK) | 0.53 | U | 0.53 | 10 |
| 2-Chloroethyl vinyl ether | 0.59 | U | 0.59 | 10 |
| 2-Chlorotoluene | 0.40 | U | 0.40 | 1.0 |
| 2-Hexanone | 0.48 | U | 0.48 | 10 |
| 4-Chlorotoluene | 0.29 | U | 0.29 | 1.0 |
| p-Isopropyltoluene | 0.43 | U | 0.43 | 1.0 |
| 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 0.99 | 10 |
| Acetone | 0.94 | U | 0.94 | 10 |
| Benzene | 0.35 | U | 0.35 | 1.0 |
| Bromobenzene | 0.35 | U | 0.35 | 1.0 |
| Bromoform | 0.56 | U | 0.56 | 1.0 |
| Bromomethane | 0.44 | U | 0.44 | 1.0 |
| Carbon disulfide | 0.38 | U | 0.38 | 1.0 |
| Carbon tetrachloride | 0.43 | U | 0.43 | 1.0 |
| Chlorobenzene | 0.25 | U | 0.25 | 1.0 |
| Dibromochloromethane | 0.43 | U | 0.43 | 1.0 |
| Chloroethane | 0.32 | U | 0.32 | 1.0 |
| Chloroform | 0.25 | U | 0.25 | 1.0 |
| Chloromethane | 0.44 | U | 0.44 | 1.0 |
| cis-1,2-Dichloroethene | 0.26 | U | 0.26 | 1.0 |
| cis-1,3-Dichloropropene | 0.46 | U | 0.46 | 1.0 |
| Dibromomethane | 0.42 | U | 0.42 | 1.0 |
| Bromochloromethane | 0.50 | U | 0.50 | 1.0 |
| Dichlorodifluoromethane | 0.32 | U | 0.32 | 1.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 1.0 |
| tert-Butyl alcohol (TBA) | 4.9 | U | 4.9 | 50 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **TB-111315**

Lab Sample ID: 240-57930-12 Date Sampled: 11/13/2015 0000
Client Matrix: Water Date Received: 11/14/2015 0945

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Prep Method: | 5030C | Prep Batch: | N/A | Lab File ID: | UXX3077.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1451 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1451 | | | | |

| Analyte | Result (ug/L) | Qualifier | MDL | RL |
|---|---|---|---|---|
| Ethylbenzene | 0.25 | U | 0.25 | 1.0 |
| Hexachlorobutadiene | 0.35 | U | 0.35 | 1.0 |
| Diisopropyl ether | 0.50 | U | 0.50 | 10 |
| Isopropylbenzene | 0.35 | U | 0.35 | 1.0 |
| Methyl tert-butyl ether | 0.20 | U | 0.20 | 1.0 |
| Methylene Chloride | 0.33 | U | 0.33 | 1.0 |
| m-Xylene & p-Xylene | 0.24 | U | 0.24 | 2.0 |
| Naphthalene | 0.45 | U | 0.45 | 1.0 |
| n-Butylbenzene | 0.31 | U | 0.31 | 1.0 |
| N-Propylbenzene | 0.40 | U | 0.40 | 1.0 |
| o-Xylene | 0.25 | U | 0.25 | 1.0 |
| sec-Butylbenzene | 0.48 | U | 0.48 | 1.0 |
| Styrene | 0.45 | U | 0.45 | 1.0 |
| Tert-amyl methyl ether | 0.30 | U | 0.30 | 5.0 |
| Ethyl t-butyl ether | 0.23 | U | 0.23 | 5.0 |
| Tetrachloroethene | 0.31 | U | 0.31 | 1.0 |
| Toluene | 0.23 | U | 0.23 | 1.0 |
| trans-1,2-Dichloroethene | 0.30 | U | 0.30 | 1.0 |
| trans-1,3-Dichloropropene | 0.56 | U | 0.56 | 1.0 |
| Trichloroethene | 0.22 | U | 0.22 | 1.0 |
| Trichlorofluoromethane | 0.49 | U | 0.49 | 1.0 |
| Vinyl acetate | 0.41 | U | 0.41 | 2.0 |
| Vinyl chloride | 0.29 | U | 0.29 | 1.0 |
| Xylenes, Total | 0.52 | U | 0.52 | 2.0 |
| tert-Butylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromoethane | 0.32 | U | 0.32 | 1.0 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 82 | | 80 - 120 |
| Dibromofluoromethane (Surr) | 81 | | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 75 | | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 83 | | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-1**

Lab Sample ID: 240-57930-1 Date Sampled: 11/13/2015 1030
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119027.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.52 g |
| Analysis Date: | 11/19/2015 1946 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 60 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.54 | U | 0.54 | 60 |
| 2-Chlorophenol | | 9.8 | U | 9.8 | 60 |
| 2-Methylnaphthalene | | 7.1 | J | 0.60 | 8.0 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.9 | U | 9.9 | 60 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 60 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 60 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.91 | U | 0.91 | 8.0 |
| Acenaphthylene | | 0.42 | U | 0.42 | 8.0 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 4.6 | J | 0.93 | 8.0 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 17 | J | 14 | 120 |
| Benzo[a]anthracene | | 12 | | 0.75 | 8.0 |
| Benzo[a]pyrene | | 12 | | 0.77 | 8.0 |
| Benzo[b]fluoranthene | | 17 | | 0.71 | 8.0 |
| Benzo[g,h,i]perylene | | 8.7 | | 0.42 | 8.0 |
| Benzo[k]fluoranthene | | 9.7 | | 0.81 | 8.0 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 23 | J B | 23 | 84 |
| Butyl benzyl phthalate | | 12 | U | 12 | 84 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 60 |
| Chrysene | | 18 | | 1.3 | 8.0 |
| Dibenz(a,h)anthracene | | 0.79 | U | 0.79 | 8.0 |
| Dibenzofuran | | 14 | J | 0.79 | 60 |
| Diethyl phthalate | | 19 | U | 19 | 84 |
| Dimethyl phthalate | | 20 | U | 20 | 84 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-1**

Lab Sample ID: 240-57930-1 Date Sampled: 11/13/2015 1030
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119027.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.52 g |
| Analysis Date: | 11/19/2015 1946 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 84 |
| Di-n-octyl phthalate | | 9.4 | U | 9.4 | 84 |
| Fluoranthene | | 27 | | 0.66 | 8.0 |
| Fluorene | | 4.8 | J | 0.63 | 8.0 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 8.0 |
| Hexachlorobutadiene | | 6.7 | U | 6.7 | 60 |
| Hexachlorocyclopentadiene | | 9.7 | U | 9.7 | 390 |
| Hexachloroethane | | 11 | U | 11 | 60 |
| Indeno[1,2,3-cd]pyrene | | 7.3 | J | 0.42 | 8.0 |
| Isophorone | | 16 | U | 16 | 60 |
| N-Nitrosodi-n-propylamine | | 7.5 | U | 7.5 | 60 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 60 |
| Naphthalene | | 7.4 | J | 0.98 | 8.0 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 20 | | 0.87 | 8.0 |
| Phenol | | 8.7 | U | 8.7 | 60 |
| Pyrene | | 23 | | 0.53 | 8.0 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 44 | | 36 - 110 |
| Phenol-d5 (Surr) | 44 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 42 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 43 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 44 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 39 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119026.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.77 g |
| Analysis Date: | 11/19/2015 1923 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 7.2 | J | 4.2 | 60 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.54 | U | 0.54 | 60 |
| 2-Chlorophenol | | 9.8 | U | 9.8 | 60 |
| 2-Methylnaphthalene | | 85 | | 0.60 | 7.9 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.9 | U | 9.9 | 60 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 60 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 60 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 10 | | 0.91 | 7.9 |
| Acenaphthylene | | 4.6 | J | 0.42 | 7.9 |
| Acetophenone | | 24 | J | 11 | 120 |
| Anthracene | | 7.4 | J | 0.93 | 7.9 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 23 | J | 14 | 120 |
| Benzo[a]anthracene | | 14 | | 0.75 | 7.9 |
| Benzo[a]pyrene | | 12 | | 0.76 | 7.9 |
| Benzo[b]fluoranthene | | 17 | | 0.70 | 7.9 |
| Benzo[g,h,i]perylene | | 10 | | 0.42 | 7.9 |
| Benzo[k]fluoranthene | | 8.9 | | 0.81 | 7.9 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 31 | J B | 23 | 83 |
| Butyl benzyl phthalate | | 12 | U | 12 | 83 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 60 |
| Chrysene | | 20 | | 1.3 | 7.9 |
| Dibenz(a,h)anthracene | | 0.79 | U | 0.79 | 7.9 |
| Dibenzofuran | | 11 | J | 0.79 | 60 |
| Diethyl phthalate | | 19 | U | 19 | 83 |
| Dimethyl phthalate | | 20 | U | 20 | 83 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119026.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.77 g |
| Analysis Date: | 11/19/2015 1923 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 83 |
| Di-n-octyl phthalate | | 9.4 | U | 9.4 | 83 |
| Fluoranthene | | 30 | | 0.66 | 7.9 |
| Fluorene | | 14 | | 0.63 | 7.9 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.9 |
| Hexachlorobutadiene | | 6.7 | U | 6.7 | 60 |
| Hexachlorocyclopentadiene | | 9.7 | U | 9.7 | 390 |
| Hexachloroethane | | 11 | U | 11 | 60 |
| Indeno[1,2,3-cd]pyrene | | 9.0 | | 0.42 | 7.9 |
| Isophorone | | 15 | U | 15 | 60 |
| N-Nitrosodi-n-propylamine | | 7.5 | U | 7.5 | 60 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 60 |
| Naphthalene | | 51 | | 0.98 | 7.9 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 35 | | 0.87 | 7.9 |
| Phenol | | 8.7 | U | 8.7 | 60 |
| Pyrene | | 28 | | 0.52 | 7.9 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 76 | | 36 - 110 |
| Phenol-d5 (Surr) | 72 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 68 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 72 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 72 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 58 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-3**

Lab Sample ID: 240-57930-3 Date Sampled: 11/13/2015 1040
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119030.D |
| Dilution: | 2.5 | | | Initial Weight/Volume: | 30.26 g |
| Analysis Date: | 11/19/2015 2057 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 85 | J | 10 | 150 |
| bis (2-chloroisopropyl) ether | | 28 | U | 28 | 290 |
| 2,4,5-Trichlorophenol | | 73 | U | 73 | 440 |
| 2,4,6-Trichlorophenol | | 26 | U | 26 | 440 |
| 2,4-Dichlorophenol | | 59 | U | 59 | 440 |
| 2,4-Dimethylphenol | | 59 | U | 59 | 440 |
| 2,4-Dinitrophenol | | 62 | U | 62 | 970 |
| 2,4-Dinitrotoluene | | 50 | U | 50 | 590 |
| 2,6-Dinitrotoluene | | 62 | U | 62 | 590 |
| 2-Chloronaphthalene | | 1.3 | U | 1.3 | 150 |
| 2-Chlorophenol | | 24 | U | 24 | 150 |
| 2-Methylnaphthalene | | 360 | | 1.5 | 20 |
| 2-Methylphenol | | 32 | U | 32 | 590 |
| 2-Nitroaniline | | 27 | U | 27 | 590 |
| 2-Nitrophenol | | 24 | U | 24 | 150 |
| 3,3'-Dichlorobenzidine | | 53 | U | 53 | 290 |
| 3-Nitroaniline | | 47 | U | 47 | 590 |
| 4,6-Dinitro-2-methylphenol | | 27 | U | 27 | 440 |
| 4-Bromophenyl phenyl ether | | 38 | U | 38 | 150 |
| 4-Chloro-3-methylphenol | | 62 | U | 62 | 440 |
| 4-Chloroaniline | | 50 | U | 50 | 440 |
| 4-Chlorophenyl phenyl ether | | 38 | U | 38 | 150 |
| 4-Nitroaniline | | 76 | U | 76 | 590 |
| 4-Nitrophenol | | 50 | U | 50 | 970 |
| Acenaphthene | | 350 | | 2.2 | 20 |
| Acenaphthylene | | 28 | | 1.0 | 20 |
| Acetophenone | | 27 | U | 27 | 290 |
| Anthracene | | 590 | | 2.3 | 20 |
| Atrazine | | 27 | U | 27 | 590 |
| Benzaldehyde | | 35 | U | 35 | 290 |
| Benzo[a]anthracene | | 960 | | 1.8 | 20 |
| Benzo[a]pyrene | | 700 | | 1.9 | 20 |
| Benzo[b]fluoranthene | | 1100 | | 1.7 | 20 |
| Benzo[g,h,i]perylene | | 240 | | 1.0 | 20 |
| Benzo[k]fluoranthene | | 440 | | 2.0 | 20 |
| Bis(2-chloroethoxy)methane | | 64 | U | 64 | 290 |
| Bis(2-chloroethyl)ether | | 5.9 | U | 5.9 | 290 |
| Bis(2-ethylhexyl) phthalate | | 56 | U | 56 | 210 |
| Butyl benzyl phthalate | | 29 | U | 29 | 210 |
| Caprolactam | | 110 | U | 110 | 970 |
| Carbazole | | 150 | | 79 | 150 |
| Chrysene | | 1000 | | 3.2 | 20 |
| Dibenz(a,h)anthracene | | 69 | | 1.9 | 20 |
| Dibenzofuran | | 330 | | 1.9 | 150 |
| Diethyl phthalate | | 47 | U | 47 | 210 |
| Dimethyl phthalate | | 50 | U | 50 | 210 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-3**

Lab Sample ID: 240-57930-3 Date Sampled: 11/13/2015 1040
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119030.D |
| Dilution: | 2.5 | | | Initial Weight/Volume: | 30.26 g |
| Analysis Date: | 11/19/2015 2057 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 44 | U | 44 | 210 |
| Di-n-octyl phthalate | | 23 | U | 23 | 210 |
| Fluoranthene | | 2500 | | 1.6 | 20 |
| Fluorene | | 350 | | 1.6 | 20 |
| Hexachlorobenzene | | 6.2 | U | 6.2 | 20 |
| Hexachlorobutadiene | | 16 | U | 16 | 150 |
| Hexachlorocyclopentadiene | | 24 | U | 24 | 970 |
| Hexachloroethane | | 26 | U | 26 | 150 |
| Indeno[1,2,3-cd]pyrene | | 250 | | 1.0 | 20 |
| Isophorone | | 38 | U | 38 | 150 |
| N-Nitrosodi-n-propylamine | | 18 | U | 18 | 150 |
| N-Nitrosodiphenylamine | | 62 | U | 62 | 150 |
| Naphthalene | | 460 | | 2.4 | 20 |
| Nitrobenzene | | 6.4 | U | 6.4 | 290 |
| Pentachlorophenol | | 27 | U | 27 | 440 |
| Phenanthrene | | 1900 | | 2.1 | 20 |
| Phenol | | 21 | U | 21 | 150 |
| Pyrene | | 1800 | | 1.3 | 20 |
| 3 & 4 Methylphenol | | 59 | U | 59 | 1200 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 76 | | 36 - 110 |
| Phenol-d5 (Surr) | 82 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 71 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 80 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 77 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 65 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-4**

Lab Sample ID: 240-57930-4 Date Sampled: 11/13/2015 1045
Client Matrix: Solid % Moisture: 16.9 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119029.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.32 g |
| Analysis Date: | 11/19/2015 2034 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 48 | J | 4.2 | 60 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.54 | U | 0.54 | 60 |
| 2-Chlorophenol | | 9.8 | U | 9.8 | 60 |
| 2-Methylnaphthalene | | 220 | | 0.60 | 7.9 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.9 | U | 9.9 | 60 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 60 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 60 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 110 | | 0.90 | 7.9 |
| Acenaphthylene | | 23 | | 0.42 | 7.9 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 160 | | 0.93 | 7.9 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 17 | J | 14 | 120 |
| Benzo[a]anthracene | | 200 | | 0.75 | 7.9 |
| Benzo[a]pyrene | | 100 | | 0.76 | 7.9 |
| Benzo[b]fluoranthene | | 240 | | 0.70 | 7.9 |
| Benzo[g,h,i]perylene | | 41 | | 0.42 | 7.9 |
| Benzo[k]fluoranthene | | 71 | | 0.81 | 7.9 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 33 | J B | 23 | 83 |
| Butyl benzyl phthalate | | 12 | U | 12 | 83 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 60 |
| Chrysene | | 260 | | 1.3 | 7.9 |
| Dibenz(a,h)anthracene | | 19 | | 0.79 | 7.9 |
| Dibenzofuran | | 170 | | 0.79 | 60 |
| Diethyl phthalate | | 19 | U | 19 | 83 |
| Dimethyl phthalate | | 20 | U | 20 | 83 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-4**

Lab Sample ID: 240-57930-4 Date Sampled: 11/13/2015 1045
Client Matrix: Solid % Moisture: 16.9 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119029.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.32 g |
| Analysis Date: | 11/19/2015 2034 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 83 |
| Di-n-octyl phthalate | | 9.4 | U | 9.4 | 83 |
| Fluoranthene | | 560 | | 0.65 | 7.9 |
| Fluorene | | 120 | | 0.63 | 7.9 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.9 |
| Hexachlorobutadiene | | 6.7 | U | 6.7 | 60 |
| Hexachlorocyclopentadiene | | 9.6 | U | 9.6 | 390 |
| Hexachloroethane | | 11 | U | 11 | 60 |
| Indeno[1,2,3-cd]pyrene | | 40 | | 0.42 | 7.9 |
| Isophorone | | 15 | U | 15 | 60 |
| N-Nitrosodi-n-propylamine | | 7.5 | U | 7.5 | 60 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 60 |
| Naphthalene | | 290 | | 0.98 | 7.9 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 550 | | 0.87 | 7.9 |
| Phenol | | 8.7 | U | 8.7 | 60 |
| Pyrene | | 420 | | 0.52 | 7.9 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 69 | | 36 - 110 |
| Phenol-d5 (Surr) | 67 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 62 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 66 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 66 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 56 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-1**

Lab Sample ID: 240-57930-5 Date Sampled: 11/13/2015 1000
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119014.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.24 g |
| Analysis Date: | 11/19/2015 1437 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 59 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 29 | U | 29 | 180 |
| 2,4,6-Trichlorophenol | | 10 | U | 10 | 180 |
| 2,4-Dichlorophenol | | 23 | U | 23 | 180 |
| 2,4-Dimethylphenol | | 23 | U | 23 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 230 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 230 |
| 2-Chloronaphthalene | | 0.53 | U | 0.53 | 59 |
| 2-Chlorophenol | | 9.6 | U | 9.6 | 59 |
| 2-Methylnaphthalene | | 0.59 | U | 0.59 | 7.8 |
| 2-Methylphenol | | 13 | U | 13 | 230 |
| 2-Nitroaniline | | 11 | U | 11 | 230 |
| 2-Nitrophenol | | 9.7 | U | 9.7 | 59 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 230 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Nitroaniline | | 30 | U | 30 | 230 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.89 | U | 0.89 | 7.8 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.8 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.91 | U | 0.91 | 7.8 |
| Atrazine | | 11 | U | 11 | 230 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.74 | U | 0.74 | 7.8 |
| Benzo[a]pyrene | | 0.75 | U | 0.75 | 7.8 |
| Benzo[b]fluoranthene | | 0.69 | U | 0.69 | 7.8 |
| Benzo[g,h,i]perylene | | 0.41 | U | 0.41 | 7.8 |
| Benzo[k]fluoranthene | | 0.80 | U | 0.80 | 7.8 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.3 | U | 2.3 | 120 |
| Bis(2-ethylhexyl) phthalate | | 25 | J B | 22 | 82 |
| Butyl benzyl phthalate | | 12 | U | 12 | 82 |
| Caprolactam | | 43 | U | 43 | 390 |
| Carbazole | | 32 | U | 32 | 59 |
| Chrysene | | 1.3 | U | 1.3 | 7.8 |
| Dibenz(a,h)anthracene | | 0.77 | U | 0.77 | 7.8 |
| Dibenzofuran | | 0.77 | U | 0.77 | 59 |
| Diethyl phthalate | | 19 | U | 19 | 82 |
| Dimethyl phthalate | | 20 | U | 20 | 82 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-1**

Lab Sample ID: 240-57930-5 Date Sampled: 11/13/2015 1000
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206908 | Lab File ID: | 51119014.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.24 g |
| Analysis Date: | 11/19/2015 1437 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 82 |
| Di-n-octyl phthalate | | 9.3 | U | 9.3 | 82 |
| Fluoranthene | | 0.64 | U | 0.64 | 7.8 |
| Fluorene | | 0.62 | U | 0.62 | 7.8 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.8 |
| Hexachlorobutadiene | | 6.6 | U | 6.6 | 59 |
| Hexachlorocyclopentadiene | | 9.5 | U | 9.5 | 390 |
| Hexachloroethane | | 11 | U | 11 | 59 |
| Indeno[1,2,3-cd]pyrene | | 0.41 | U | 0.41 | 7.8 |
| Isophorone | | 15 | U | 15 | 59 |
| N-Nitrosodi-n-propylamine | | 7.4 | U | 7.4 | 59 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 59 |
| Naphthalene | | 0.96 | U | 0.96 | 7.8 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.86 | U | 0.86 | 7.8 |
| Phenol | | 8.6 | U | 8.6 | 59 |
| Pyrene | | 0.52 | U | 0.52 | 7.8 |
| 3 & 4 Methylphenol | | 23 | U | 23 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 86 | | 36 - 110 |
| Phenol-d5 (Surr) | 80 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 70 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 78 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 79 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 56 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-2**

Lab Sample ID: 240-57930-6 Date Sampled: 11/13/2015 1005
Client Matrix: Solid % Moisture: 15.5 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119017.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.05 g |
| Analysis Date: | 11/19/2015 1548 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 59 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.53 | U | 0.53 | 59 |
| 2-Chlorophenol | | 9.7 | U | 9.7 | 59 |
| 2-Methylnaphthalene | | 0.59 | U | 0.59 | 7.9 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.8 | U | 9.8 | 59 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.90 | U | 0.90 | 7.9 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.9 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.92 | U | 0.92 | 7.9 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.74 | U | 0.74 | 7.9 |
| Benzo[a]pyrene | | 0.76 | U | 0.76 | 7.9 |
| Benzo[b]fluoranthene | | 0.70 | U | 0.70 | 7.9 |
| Benzo[g,h,i]perylene | | 0.41 | U | 0.41 | 7.9 |
| Benzo[k]fluoranthene | | 0.80 | U | 0.80 | 7.9 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 38 | J B | 22 | 83 |
| Butyl benzyl phthalate | | 12 | U | 12 | 83 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 59 |
| Chrysene | | 1.3 | U | 1.3 | 7.9 |
| Dibenz(a,h)anthracene | | 0.78 | U | 0.78 | 7.9 |
| Dibenzofuran | | 0.78 | U | 0.78 | 59 |
| Diethyl phthalate | | 19 | U | 19 | 83 |
| Dimethyl phthalate | | 20 | U | 20 | 83 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-2**

Lab Sample ID: 240-57930-6 Date Sampled: 11/13/2015 1005
Client Matrix: Solid % Moisture: 15.5 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119017.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.05 g |
| Analysis Date: | 11/19/2015 1548 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 83 |
| Di-n-octyl phthalate | | 9.3 | U | 9.3 | 83 |
| Fluoranthene | | 0.65 | U | 0.65 | 7.9 |
| Fluorene | | 0.63 | U | 0.63 | 7.9 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.9 |
| Hexachlorobutadiene | | 6.6 | U | 6.6 | 59 |
| Hexachlorocyclopentadiene | | 9.6 | U | 9.6 | 390 |
| Hexachloroethane | | 11 | U | 11 | 59 |
| Indeno[1,2,3-cd]pyrene | | 0.41 | U | 0.41 | 7.9 |
| Isophorone | | 15 | U | 15 | 59 |
| N-Nitrosodi-n-propylamine | | 7.4 | U | 7.4 | 59 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 59 |
| Naphthalene | | 0.97 | U | 0.97 | 7.9 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.86 | U | 0.86 | 7.9 |
| Phenol | | 8.6 | U | 8.6 | 59 |
| Pyrene | | 0.52 | U | 0.52 | 7.9 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 68 | | 36 - 110 |
| Phenol-d5 (Surr) | 68 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 64 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 68 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 67 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 55 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-3**

Lab Sample ID: 240-57930-7 Date Sampled: 11/13/2015 1010
Client Matrix: Solid % Moisture: 19.0 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119018.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.21 g |
| Analysis Date: | 11/19/2015 1612 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.3 | U | 4.3 | 61 |
| bis (2-chloroisopropyl) ether | | 12 | U | 12 | 120 |
| 2,4,5-Trichlorophenol | | 31 | U | 31 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 25 | U | 25 | 180 |
| 2,4-Dimethylphenol | | 25 | U | 25 | 180 |
| 2,4-Dinitrophenol | | 26 | U | 26 | 400 |
| 2,4-Dinitrotoluene | | 21 | U | 21 | 250 |
| 2,6-Dinitrotoluene | | 26 | U | 26 | 250 |
| 2-Chloronaphthalene | | 0.55 | U | 0.55 | 61 |
| 2-Chlorophenol | | 10 | U | 10 | 61 |
| 2-Methylnaphthalene | | 0.61 | U | 0.61 | 8.2 |
| 2-Methylphenol | | 13 | U | 13 | 250 |
| 2-Nitroaniline | | 11 | U | 11 | 250 |
| 2-Nitrophenol | | 10 | U | 10 | 61 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 20 | U | 20 | 250 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Chloro-3-methylphenol | | 26 | U | 26 | 180 |
| 4-Chloroaniline | | 21 | U | 21 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Nitroaniline | | 32 | U | 32 | 250 |
| 4-Nitrophenol | | 21 | U | 21 | 400 |
| Acenaphthene | | 0.93 | U | 0.93 | 8.2 |
| Acenaphthylene | | 0.43 | U | 0.43 | 8.2 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.96 | U | 0.96 | 8.2 |
| Atrazine | | 11 | U | 11 | 250 |
| Benzaldehyde | | 45 | J | 15 | 120 |
| Benzo[a]anthracene | | 0.77 | U | 0.77 | 8.2 |
| Benzo[a]pyrene | | 0.78 | U | 0.78 | 8.2 |
| Benzo[b]fluoranthene | | 0.72 | U | 0.72 | 8.2 |
| Benzo[g,h,i]perylene | | 0.43 | U | 0.43 | 8.2 |
| Benzo[k]fluoranthene | | 0.83 | U | 0.83 | 8.2 |
| Bis(2-chloroethoxy)methane | | 27 | U | 27 | 120 |
| Bis(2-chloroethyl)ether | | 2.5 | U | 2.5 | 120 |
| Bis(2-ethylhexyl) phthalate | | 55 | J B | 23 | 86 |
| Butyl benzyl phthalate | | 12 | U | 12 | 86 |
| Caprolactam | | 45 | U | 45 | 400 |
| Carbazole | | 33 | U | 33 | 61 |
| Chrysene | | 1.3 | U | 1.3 | 8.2 |
| Dibenz(a,h)anthracene | | 0.81 | U | 0.81 | 8.2 |
| Dibenzofuran | | 0.81 | U | 0.81 | 61 |
| Diethyl phthalate | | 20 | U | 20 | 86 |
| Dimethyl phthalate | | 21 | U | 21 | 86 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-3**

Lab Sample ID: 240-57930-7 Date Sampled: 11/13/2015 1010
Client Matrix: Solid % Moisture: 19.0 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119018.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.21 g |
| Analysis Date: | 11/19/2015 1612 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 86 |
| Di-n-octyl phthalate | | 9.7 | U | 9.7 | 86 |
| Fluoranthene | | 0.67 | U | 0.67 | 8.2 |
| Fluorene | | 0.65 | U | 0.65 | 8.2 |
| Hexachlorobenzene | | 2.6 | U | 2.6 | 8.2 |
| Hexachlorobutadiene | | 6.9 | U | 6.9 | 61 |
| Hexachlorocyclopentadiene | | 9.9 | U | 9.9 | 400 |
| Hexachloroethane | | 11 | U | 11 | 61 |
| Indeno[1,2,3-cd]pyrene | | 0.43 | U | 0.43 | 8.2 |
| Isophorone | | 16 | U | 16 | 61 |
| N-Nitrosodi-n-propylamine | | 7.7 | U | 7.7 | 61 |
| N-Nitrosodiphenylamine | | 26 | U | 26 | 61 |
| Naphthalene | | 1.0 | U | 1.0 | 8.2 |
| Nitrobenzene | | 2.7 | U | 2.7 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.90 | U | 0.90 | 8.2 |
| Phenol | | 9.0 | U | 9.0 | 61 |
| Pyrene | | 0.54 | U | 0.54 | 8.2 |
| 3 & 4 Methylphenol | | 25 | U | 25 | 490 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 66 | | 36 - 110 |
| Phenol-d5 (Surr) | 67 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 59 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 66 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 63 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 40 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-4**

Lab Sample ID: 240-57930-8 Date Sampled: 11/13/2015 1015
Client Matrix: Solid % Moisture: 16.1 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119019.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.80 g |
| Analysis Date: | 11/19/2015 1636 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 60 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 400 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.54 | U | 0.54 | 60 |
| 2-Chlorophenol | | 9.8 | U | 9.8 | 60 |
| 2-Methylnaphthalene | | 0.60 | U | 0.60 | 8.0 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 10 | U | 10 | 60 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 60 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 60 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 400 |
| Acenaphthene | | 0.91 | U | 0.91 | 8.0 |
| Acenaphthylene | | 0.42 | U | 0.42 | 8.0 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.94 | U | 0.94 | 8.0 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.76 | U | 0.76 | 8.0 |
| Benzo[a]pyrene | | 0.77 | U | 0.77 | 8.0 |
| Benzo[b]fluoranthene | | 0.71 | U | 0.71 | 8.0 |
| Benzo[g,h,i]perylene | | 0.42 | U | 0.42 | 8.0 |
| Benzo[k]fluoranthene | | 0.82 | U | 0.82 | 8.0 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 30 | J B | 23 | 84 |
| Butyl benzyl phthalate | | 12 | U | 12 | 84 |
| Caprolactam | | 44 | U | 44 | 400 |
| Carbazole | | 32 | U | 32 | 60 |
| Chrysene | | 1.3 | U | 1.3 | 8.0 |
| Dibenz(a,h)anthracene | | 0.79 | U | 0.79 | 8.0 |
| Dibenzofuran | | 0.79 | U | 0.79 | 60 |
| Diethyl phthalate | | 19 | U | 19 | 84 |
| Dimethyl phthalate | | 20 | U | 20 | 84 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-4**

Lab Sample ID: 240-57930-8 Date Sampled: 11/13/2015 1015
Client Matrix: Solid % Moisture: 16.1 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119019.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.80 g |
| Analysis Date: | 11/19/2015 1636 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 84 |
| Di-n-octyl phthalate | | 9.5 | U | 9.5 | 84 |
| Fluoranthene | | 0.66 | U | 0.66 | 8.0 |
| Fluorene | | 0.64 | U | 0.64 | 8.0 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 8.0 |
| Hexachlorobutadiene | | 6.7 | U | 6.7 | 60 |
| Hexachlorocyclopentadiene | | 9.7 | U | 9.7 | 400 |
| Hexachloroethane | | 11 | U | 11 | 60 |
| Indeno[1,2,3-cd]pyrene | | 0.42 | U | 0.42 | 8.0 |
| Isophorone | | 16 | U | 16 | 60 |
| N-Nitrosodi-n-propylamine | | 7.6 | U | 7.6 | 60 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 60 |
| Naphthalene | | 0.98 | U | 0.98 | 8.0 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.88 | U | 0.88 | 8.0 |
| Phenol | | 8.8 | U | 8.8 | 60 |
| Pyrene | | 0.53 | U | 0.53 | 8.0 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 74 | | 36 - 110 |
| Phenol-d5 (Surr) | 75 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 66 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 73 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 71 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 54 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-5**

Lab Sample ID: 240-57930-9 Date Sampled: 11/13/2015 1020
Client Matrix: Solid % Moisture: 16.7 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119020.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.71 g |
| Analysis Date: | 11/19/2015 1700 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.2 | U | 4.2 | 61 |
| bis (2-chloroisopropyl) ether | | 12 | U | 12 | 120 |
| 2,4,5-Trichlorophenol | | 30 | U | 30 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 400 |
| 2,4-Dinitrotoluene | | 21 | U | 21 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.55 | U | 0.55 | 61 |
| 2-Chlorophenol | | 9.9 | U | 9.9 | 61 |
| 2-Methylnaphthalene | | 0.61 | U | 0.61 | 8.1 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 10 | U | 10 | 61 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 21 | U | 21 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Nitroaniline | | 32 | U | 32 | 240 |
| 4-Nitrophenol | | 21 | U | 21 | 400 |
| Acenaphthene | | 0.92 | U | 0.92 | 8.1 |
| Acenaphthylene | | 0.42 | U | 0.42 | 8.1 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.95 | U | 0.95 | 8.1 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 15 | U | 15 | 120 |
| Benzo[a]anthracene | | 0.76 | U | 0.76 | 8.1 |
| Benzo[a]pyrene | | 0.78 | U | 0.78 | 8.1 |
| Benzo[b]fluoranthene | | 0.72 | U | 0.72 | 8.1 |
| Benzo[g,h,i]perylene | | 0.42 | U | 0.42 | 8.1 |
| Benzo[k]fluoranthene | | 0.82 | U | 0.82 | 8.1 |
| Bis(2-chloroethoxy)methane | | 27 | U | 27 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 62 | J B | 23 | 85 |
| Butyl benzyl phthalate | | 12 | U | 12 | 85 |
| Caprolactam | | 45 | U | 45 | 400 |
| Carbazole | | 33 | U | 33 | 61 |
| Chrysene | | 1.3 | U | 1.3 | 8.1 |
| Dibenz(a,h)anthracene | | 0.80 | U | 0.80 | 8.1 |
| Dibenzofuran | | 0.80 | U | 0.80 | 61 |
| Diethyl phthalate | | 19 | U | 19 | 85 |
| Dimethyl phthalate | | 21 | U | 21 | 85 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-5**

Lab Sample ID: 240-57930-9 Date Sampled: 11/13/2015 1020
Client Matrix: Solid % Moisture: 16.7 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119020.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 29.71 g |
| Analysis Date: | 11/19/2015 1700 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 85 |
| Di-n-octyl phthalate | | 9.6 | U | 9.6 | 85 |
| Fluoranthene | | 0.67 | U | 0.67 | 8.1 |
| Fluorene | | 0.64 | U | 0.64 | 8.1 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 8.1 |
| Hexachlorobutadiene | | 6.8 | U | 6.8 | 61 |
| Hexachlorocyclopentadiene | | 9.8 | U | 9.8 | 400 |
| Hexachloroethane | | 11 | U | 11 | 61 |
| Indeno[1,2,3-cd]pyrene | | 0.42 | U | 0.42 | 8.1 |
| Isophorone | | 16 | U | 16 | 61 |
| N-Nitrosodi-n-propylamine | | 7.6 | U | 7.6 | 61 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 61 |
| Naphthalene | | 0.99 | U | 0.99 | 8.1 |
| Nitrobenzene | | 2.7 | U | 2.7 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.88 | U | 0.88 | 8.1 |
| Phenol | | 8.8 | U | 8.8 | 61 |
| Pyrene | | 0.53 | U | 0.53 | 8.1 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 480 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 73 | | 36 - 110 |
| Phenol-d5 (Surr) | 69 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 60 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 68 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 67 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 41 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-6**

Lab Sample ID: 240-57930-10 Date Sampled: 11/13/2015 1025
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119021.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.14 g |
| Analysis Date: | 11/19/2015 1723 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 59 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 29 | U | 29 | 180 |
| 2,4,6-Trichlorophenol | | 10 | U | 10 | 180 |
| 2,4-Dichlorophenol | | 23 | U | 23 | 180 |
| 2,4-Dimethylphenol | | 23 | U | 23 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 230 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 230 |
| 2-Chloronaphthalene | | 0.53 | U | 0.53 | 59 |
| 2-Chlorophenol | | 9.6 | U | 9.6 | 59 |
| 2-Methylnaphthalene | | 0.59 | U | 0.59 | 7.8 |
| 2-Methylphenol | | 13 | U | 13 | 230 |
| 2-Nitroaniline | | 11 | U | 11 | 230 |
| 2-Nitrophenol | | 9.7 | U | 9.7 | 59 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 230 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Nitroaniline | | 31 | U | 31 | 230 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.89 | U | 0.89 | 7.8 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.8 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.92 | U | 0.92 | 7.8 |
| Atrazine | | 11 | U | 11 | 230 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 0.74 | U | 0.74 | 7.8 |
| Benzo[a]pyrene | | 0.75 | U | 0.75 | 7.8 |
| Benzo[b]fluoranthene | | 0.69 | U | 0.69 | 7.8 |
| Benzo[g,h,i]perylene | | 0.41 | U | 0.41 | 7.8 |
| Benzo[k]fluoranthene | | 0.80 | U | 0.80 | 7.8 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.3 | U | 2.3 | 120 |
| Bis(2-ethylhexyl) phthalate | | 31 | J B | 22 | 82 |
| Butyl benzyl phthalate | | 12 | U | 12 | 82 |
| Caprolactam | | 43 | U | 43 | 390 |
| Carbazole | | 32 | U | 32 | 59 |
| Chrysene | | 1.3 | U | 1.3 | 7.8 |
| Dibenz(a,h)anthracene | | 0.77 | U | 0.77 | 7.8 |
| Dibenzofuran | | 0.77 | U | 0.77 | 59 |
| Diethyl phthalate | | 19 | U | 19 | 82 |
| Dimethyl phthalate | | 20 | U | 20 | 82 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-6**

Lab Sample ID: 240-57930-10 Date Sampled: 11/13/2015 1025
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119021.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.14 g |
| Analysis Date: | 11/19/2015 1723 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 82 |
| Di-n-octyl phthalate | | 9.3 | U | 9.3 | 82 |
| Fluoranthene | | 0.65 | U | 0.65 | 7.8 |
| Fluorene | | 0.62 | U | 0.62 | 7.8 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.8 |
| Hexachlorobutadiene | | 6.6 | U | 6.6 | 59 |
| Hexachlorocyclopentadiene | | 9.5 | U | 9.5 | 390 |
| Hexachloroethane | | 11 | U | 11 | 59 |
| Indeno[1,2,3-cd]pyrene | | 0.41 | U | 0.41 | 7.8 |
| Isophorone | | 15 | U | 15 | 59 |
| N-Nitrosodi-n-propylamine | | 7.4 | U | 7.4 | 59 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 59 |
| Naphthalene | | 0.96 | U | 0.96 | 7.8 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 0.86 | U | 0.86 | 7.8 |
| Phenol | | 8.6 | U | 8.6 | 59 |
| Pyrene | | 0.52 | U | 0.52 | 7.8 |
| 3 & 4 Methylphenol | | 23 | U | 23 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 70 | | 36 - 110 |
| Phenol-d5 (Surr) | 69 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 61 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 68 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 66 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 53 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-DUP**

Lab Sample ID: 240-57930-11 Date Sampled: 11/13/2015 0000
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119028.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.46 g |
| Analysis Date: | 11/19/2015 2010 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 58 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 29 | U | 29 | 180 |
| 2,4,6-Trichlorophenol | | 10 | U | 10 | 180 |
| 2,4-Dichlorophenol | | 23 | U | 23 | 180 |
| 2,4-Dimethylphenol | | 23 | U | 23 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 230 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 230 |
| 2-Chloronaphthalene | | 0.53 | U | 0.53 | 58 |
| 2-Chlorophenol | | 9.6 | U | 9.6 | 58 |
| 2-Methylnaphthalene | | 7.8 | | 0.58 | 7.8 |
| 2-Methylphenol | | 13 | U | 13 | 230 |
| 2-Nitroaniline | | 11 | U | 11 | 230 |
| 2-Nitrophenol | | 9.7 | U | 9.7 | 58 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 230 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 58 |
| 4-Nitroaniline | | 30 | U | 30 | 230 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 0.89 | U | 0.89 | 7.8 |
| Acenaphthylene | | 4.0 | J | 0.41 | 7.8 |
| Acetophenone | | 17 | J | 11 | 120 |
| Anthracene | | 5.4 | J | 0.91 | 7.8 |
| Atrazine | | 11 | U | 11 | 230 |
| Benzaldehyde | | 25 | J | 14 | 120 |
| Benzo[a]anthracene | | 16 | | 0.74 | 7.8 |
| Benzo[a]pyrene | | 17 | | 0.75 | 7.8 |
| Benzo[b]fluoranthene | | 30 | | 0.69 | 7.8 |
| Benzo[g,h,i]perylene | | 10 | | 0.41 | 7.8 |
| Benzo[k]fluoranthene | | 9.9 | | 0.79 | 7.8 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.3 | U | 2.3 | 120 |
| Bis(2-ethylhexyl) phthalate | | 22 | U | 22 | 82 |
| Butyl benzyl phthalate | | 12 | U | 12 | 82 |
| Caprolactam | | 43 | U | 43 | 390 |
| Carbazole | | 32 | U | 32 | 58 |
| Chrysene | | 22 | | 1.3 | 7.8 |
| Dibenz(a,h)anthracene | | 0.77 | U | 0.77 | 7.8 |
| Dibenzofuran | | 16 | J | 0.77 | 58 |
| Diethyl phthalate | | 19 | U | 19 | 82 |
| Dimethyl phthalate | | 20 | U | 20 | 82 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-DUP**

Lab Sample ID: 240-57930-11 Date Sampled: 11/13/2015 0000
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/14/2015 0945

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Prep Method: | 3540C | Prep Batch: | 240-206887 | Lab File ID: | 51119028.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.46 g |
| Analysis Date: | 11/19/2015 2010 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 18 | U | 18 | 82 |
| Di-n-octyl phthalate | | 9.2 | U | 9.2 | 82 |
| Fluoranthene | | 37 | | 0.64 | 7.8 |
| Fluorene | | 6.1 | J | 0.62 | 7.8 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.8 |
| Hexachlorobutadiene | | 6.5 | U | 6.5 | 58 |
| Hexachlorocyclopentadiene | | 9.5 | U | 9.5 | 390 |
| Hexachloroethane | | 11 | U | 11 | 58 |
| Indeno[1,2,3-cd]pyrene | | 9.2 | | 0.41 | 7.8 |
| Isophorone | | 15 | U | 15 | 58 |
| N-Nitrosodi-n-propylamine | | 7.4 | U | 7.4 | 58 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 58 |
| Naphthalene | | 12 | | 0.96 | 7.8 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 27 | | 0.85 | 7.8 |
| Phenol | | 8.5 | U | 8.5 | 58 |
| Pyrene | | 30 | | 0.51 | 7.8 |
| 3 & 4 Methylphenol | | 23 | U | 23 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 66 | | 36 - 110 |
| Phenol-d5 (Surr) | 68 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 65 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 68 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 66 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 62 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-1**

Lab Sample ID: 240-57930-1 Date Sampled: 11/13/2015 1030
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-206878 | Initial Weight/Volume: | 0.55 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/16/2015 1118 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 3500 | | 490 | 1100 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 69 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207037 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-207078 | Initial Weight/Volume: | 0.48 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1039 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/17/2015 0748 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 5500 | B | 570 | 1200 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 78 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-3**

Lab Sample ID: 240-57930-3 Date Sampled: 11/13/2015 1040
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207246 | Instrument ID: | AFID |
| Prep Method: | 5030A | Prep Batch: | 240-207274 | Initial Weight/Volume: | 5.02 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/18/2015 1704 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/18/2015 0657 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 110 | J | 54 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 75 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VS-UST-P-4**

Lab Sample ID: 240-57930-4 Date Sampled: 11/13/2015 1045
Client Matrix: Solid % Moisture: 16.9 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-206878 | Initial Weight/Volume: | 5.01 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/16/2015 1301 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 67 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-1**

Lab Sample ID: 240-57930-5 Date Sampled: 11/13/2015 1000
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-206878 | Initial Weight/Volume: | 5.05 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/16/2015 1336 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 54 | U | 54 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 72 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID: F-VB-UST-P-2**

Lab Sample ID: 240-57930-6 Date Sampled: 11/13/2015 1005
Client Matrix: Solid % Moisture: 15.5 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-206878 | Initial Weight/Volume: | 4.96 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/16/2015 1412 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 64 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-3**

Lab Sample ID: 240-57930-7 Date Sampled: 11/13/2015 1010
Client Matrix: Solid % Moisture: 19.0 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-206878 | Initial Weight/Volume: | 5.08 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/16/2015 1519 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 60 | J | 56 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 72 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-4**

Lab Sample ID: 240-57930-8 Date Sampled: 11/13/2015 1015
Client Matrix: Solid % Moisture: 16.1 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-206878 | Initial Weight/Volume: | 4.95 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/16/2015 1556 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 53 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-5**

Lab Sample ID: 240-57930-9 Date Sampled: 11/13/2015 1020
Client Matrix: Solid % Moisture: 16.7 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-206878 | Initial Weight/Volume: | 4.98 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/16/2015 1630 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 51 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-6**

Lab Sample ID: 240-57930-10 Date Sampled: 11/13/2015 1025
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-206878 | Initial Weight/Volume: | 4.96 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/16/2015 1705 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 72 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-DUP**

Lab Sample ID: 240-57930-11 Date Sampled: 11/13/2015 0000
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-206878 | Initial Weight/Volume: | 0.50 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/16/2015 1740 | | | Injection Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 1800 | | 540 | 1200 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 70 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-1**

Lab Sample ID: 240-57930-1 Date Sampled: 11/13/2015 1030
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207234 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-206884 | Initial Weight/Volume: | 30.28 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/18/2015 0732 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0833 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 80 | | 7.5 | 19 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 82 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207234 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-206884 | Initial Weight/Volume: | 29.96 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/18/2015 0804 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0833 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 130 | | 7.6 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 95 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-3**

Lab Sample ID: 240-57930-3 Date Sampled: 11/13/2015 1040
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207234 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-206884 | Initial Weight/Volume: | 30.34 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/18/2015 0835 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0833 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 42 | | 7.5 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 75 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-4**

Lab Sample ID: 240-57930-4 Date Sampled: 11/13/2015 1045
Client Matrix: Solid % Moisture: 16.9 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207234 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-206884 | Initial Weight/Volume: | 30.19 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/18/2015 1010 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0833 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 120 | | 7.7 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 82 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-1**

Lab Sample ID: 240-57930-5 Date Sampled: 11/13/2015 1000
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207234 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-206884 | Initial Weight/Volume: | 29.96 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/18/2015 1047 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0833 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.6 | U | 7.6 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 76 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-2**

Lab Sample ID: 240-57930-6 Date Sampled: 11/13/2015 1005
Client Matrix: Solid % Moisture: 15.5 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207234 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-206884 | Initial Weight/Volume: | 29.88 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/18/2015 1121 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0833 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.6 | U | 7.6 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 79 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-3**

Lab Sample ID: 240-57930-7 Date Sampled: 11/13/2015 1010
Client Matrix: Solid % Moisture: 19.0 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207234 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-206884 | Initial Weight/Volume: | 30.03 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/18/2015 1232 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0833 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.9 | U | 7.9 | 21 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 69 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-4**

Lab Sample ID: 240-57930-8 Date Sampled: 11/13/2015 1015
Client Matrix: Solid % Moisture: 16.1 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207465 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-207121 | Initial Weight/Volume: | 30.20 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/19/2015 0232 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/17/2015 1002 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.6 | U | 7.6 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 90 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-5**

Lab Sample ID: 240-57930-9 Date Sampled: 11/13/2015 1020
Client Matrix: Solid % Moisture: 16.7 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207465 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-207121 | Initial Weight/Volume: | 29.52 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/19/2015 0304 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/17/2015 1002 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.8 | U | 7.8 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 81 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-6**

Lab Sample ID: 240-57930-10 Date Sampled: 11/13/2015 1025
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207465 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-207121 | Initial Weight/Volume: | 30.04 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/19/2015 0336 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/17/2015 1002 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 8.0 | J | 7.5 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 83 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-DUP**

Lab Sample ID: 240-57930-11 Date Sampled: 11/13/2015 0000
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/14/2015 0945

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-207465 | Instrument ID: | A2HP5R |
| Prep Method: | 3540C | Prep Batch: | 240-207121 | Initial Weight/Volume: | 30.26 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 11/19/2015 0407 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/17/2015 1002 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 51 | | 7.5 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 79 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-1**

Lab Sample ID: 240-57930-1 Date Sampled: 11/13/2015 1030
Client Matrix: Solid % Moisture: 15.0 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 29.62 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/18/2015 2351 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 24 | U | 24 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.5 | U | 9.5 | 39 |
| Aroclor-1254 | | 17 | U | 17 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 83 | | 13 - 134 |
| DCB Decachlorobiphenyl | 72 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 30.44 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/19/2015 0009 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 38 |
| Aroclor-1221 | | 19 | U | 19 | 38 |
| Aroclor-1232 | | 23 | U | 23 | 38 |
| Aroclor-1242 | | 13 | U | 13 | 38 |
| Aroclor-1248 | | 9.3 | U | 9.3 | 38 |
| Aroclor-1254 | | 16 | U | 16 | 38 |
| Aroclor-1260 | | 10 | U | 10 | 38 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 61 | | 13 - 134 |
| DCB Decachlorobiphenyl | 70 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-3**

Lab Sample ID: 240-57930-3 Date Sampled: 11/13/2015 1040
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 29.65 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/19/2015 0027 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 24 | U | 24 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.6 | U | 9.6 | 39 |
| Aroclor-1254 | | 17 | U | 17 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 67 | | 13 - 134 |
| DCB Decachlorobiphenyl | 90 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-4**

Lab Sample ID: 240-57930-4 Date Sampled: 11/13/2015 1045
Client Matrix: Solid % Moisture: 16.9 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 30.41 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/19/2015 0045 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 24 | U | 24 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.5 | U | 9.5 | 39 |
| Aroclor-1254 | | 17 | U | 17 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 78 | | 13 - 134 |
| DCB Decachlorobiphenyl | 141 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-1**

Lab Sample ID: 240-57930-5 Date Sampled: 11/13/2015 1000
Client Matrix: Solid % Moisture: 15.4 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 30.45 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/19/2015 0158 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 38 |
| Aroclor-1221 | | 19 | U | 19 | 38 |
| Aroclor-1232 | | 23 | U | 23 | 38 |
| Aroclor-1242 | | 13 | U | 13 | 38 |
| Aroclor-1248 | | 9.3 | U | 9.3 | 38 |
| Aroclor-1254 | | 16 | U | 16 | 38 |
| Aroclor-1260 | | 10 | U | 10 | 38 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 83 | | 13 - 134 |
| DCB Decachlorobiphenyl | 82 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-2**

Lab Sample ID: 240-57930-6 Date Sampled: 11/13/2015 1005
Client Matrix: Solid % Moisture: 15.5 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 29.88 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/18/2015 2202 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 24 | U | 24 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.5 | U | 9.5 | 39 |
| Aroclor-1254 | | 17 | U | 17 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 70 | | 13 - 134 |
| DCB Decachlorobiphenyl | 66 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-3**

Lab Sample ID: 240-57930-7 Date Sampled: 11/13/2015 1010
Client Matrix: Solid % Moisture: 19.0 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 30.01 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/18/2015 2220 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 15 | U | 15 | 41 |
| Aroclor-1221 | | 20 | U | 20 | 41 |
| Aroclor-1232 | | 25 | U | 25 | 41 |
| Aroclor-1242 | | 14 | U | 14 | 41 |
| Aroclor-1248 | | 9.9 | U | 9.9 | 41 |
| Aroclor-1254 | | 17 | U | 17 | 41 |
| Aroclor-1260 | | 11 | U | 11 | 41 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 67 | | 13 - 134 |
| DCB Decachlorobiphenyl | 66 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-4**

Lab Sample ID: 240-57930-8 Date Sampled: 11/13/2015 1015
Client Matrix: Solid % Moisture: 16.1 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 29.74 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/18/2015 2238 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 40 |
| Aroclor-1221 | | 19 | U | 19 | 40 |
| Aroclor-1232 | | 24 | U | 24 | 40 |
| Aroclor-1242 | | 13 | U | 13 | 40 |
| Aroclor-1248 | | 9.6 | U | 9.6 | 40 |
| Aroclor-1254 | | 17 | U | 17 | 40 |
| Aroclor-1260 | | 11 | U | 11 | 40 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 83 | | 13 - 134 |
| DCB Decachlorobiphenyl | 79 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-5**

Lab Sample ID: 240-57930-9 Date Sampled: 11/13/2015 1020
Client Matrix: Solid % Moisture: 16.7 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 29.69 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/18/2015 2256 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 15 | U | 15 | 40 |
| Aroclor-1221 | | 19 | U | 19 | 40 |
| Aroclor-1232 | | 24 | U | 24 | 40 |
| Aroclor-1242 | | 13 | U | 13 | 40 |
| Aroclor-1248 | | 9.7 | U | 9.7 | 40 |
| Aroclor-1254 | | 17 | U | 17 | 40 |
| Aroclor-1260 | | 11 | U | 11 | 40 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 80 | | 13 - 134 |
| DCB Decachlorobiphenyl | 71 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VB-UST-P-6**

Lab Sample ID: 240-57930-10 Date Sampled: 11/13/2015 1025
Client Matrix: Solid % Moisture: 15.2 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 29.68 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/18/2015 2315 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 24 | U | 24 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.5 | U | 9.5 | 39 |
| Aroclor-1254 | | 17 | U | 17 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 72 | | 13 - 134 |
| DCB Decachlorobiphenyl | 71 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Client Sample ID:** **F-VS-UST-P-DUP**

Lab Sample ID: 240-57930-11 Date Sampled: 11/13/2015 0000
Client Matrix: Solid % Moisture: 15.6 Date Received: 11/14/2015 0945

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-206917 | Initial Weight/Volume: | 29.83 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 11/18/2015 2333 | | | Injection Volume: | 1 uL |
| Prep Date: | 11/16/2015 0932 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 39 |
| Aroclor-1221 | | 19 | U | 19 | 39 |
| Aroclor-1232 | | 24 | U | 24 | 39 |
| Aroclor-1242 | | 13 | U | 13 | 39 |
| Aroclor-1248 | | 9.5 | U | 9.5 | 39 |
| Aroclor-1254 | | 17 | U | 17 | 39 |
| Aroclor-1260 | | 11 | U | 11 | 39 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 82 | | 13 - 134 |
| DCB Decachlorobiphenyl | 78 | | 10 - 155 |

**General Chemistry**

**Client Sample ID:** **F-VS-UST-P-1**

Lab Sample ID: 240-57930-1 Date Sampled: 11/13/2015 1030
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 85 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 15 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VS-UST-P-2**

Lab Sample ID: 240-57930-2 Date Sampled: 11/13/2015 1035
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 85 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 15 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VS-UST-P-3**

Lab Sample ID: 240-57930-3 Date Sampled: 11/13/2015 1040
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 85 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 15 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VS-UST-P-4**

Lab Sample ID: 240-57930-4 Date Sampled: 11/13/2015 1045
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 83 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 17 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-P-1**

Lab Sample ID: 240-57930-5 Date Sampled: 11/13/2015 1000
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 85 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 15 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-P-2**

Lab Sample ID: 240-57930-6 Date Sampled: 11/13/2015 1005
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 84 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 16 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-P-3**

Lab Sample ID: 240-57930-7 Date Sampled: 11/13/2015 1010
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 81 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 19 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-P-4**

Lab Sample ID: 240-57930-8 Date Sampled: 11/13/2015 1015
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 84 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 16 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-P-5**

Lab Sample ID: 240-57930-9 Date Sampled: 11/13/2015 1020
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 83 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 17 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-P-6**

Lab Sample ID: 240-57930-10 Date Sampled: 11/13/2015 1025
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 85 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 15 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VS-UST-P-DUP**

Lab Sample ID: 240-57930-11 Date Sampled: 11/13/2015 0000
Client Matrix: Solid Date Received: 11/14/2015 0945

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 84 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |
| Percent Moisture | 16 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-207016 | Analysis Date: 11/16/2015 1541 | | | | | | DryWt Corrected: N |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Surrogate Recovery Report

**8260C Volatile Organic Compounds by GC/MS**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | DBFM %Rec | DCA %Rec | TOL %Rec | BFB %Rec |
|---|---|---|---|---|---|
| 240-57930-1 | F-VS-UST-P-1 | 100 | 104 | 104 | 97 |
| 240-57930-2 | F-VS-UST-P-2 | 92 | 98 | 94 | 91 |
| 240-57930-2 | F-VS-UST-P-2 | 92 | 95 | 103 | 105 |
| 240-57930-2 | F-VS-UST-P-2 | 94 | 88 | 110 | 93 |
| 240-57930-3 | F-VS-UST-P-3 | 91 | 91 | 94 | 89 |
| 240-57930-4 | F-VS-UST-P-4 | 87 | 85 | 111 | 102* |
| 240-57930-5 | F-VB-UST-P-1 | 89 | 90 | 89 | 78 |
| 240-57930-6 | F-VB-UST-P-2 | 95 | 100 | 89 | 79 |
| 240-57930-7 | F-VB-UST-P-3 | 99 | 105 | 94 | 82 |
| 240-57930-8 | F-VB-UST-P-4 | 103 | 113 | 96 | 89 |
| 240-57930-9 | F-VB-UST-P-5 | 86 | 89 | 84 | 76 |
| 240-57930-10 | F-VB-UST-P-6 | 97 | 105 | 93 | 86 |
| 240-57930-11 | F-VS-UST-P-DUP | 89 | 94 | 90 | 83 |
| MB 240-207343/6 | | 91 | 95 | 87 | 76 |
| MB 240-207560/6 | | 98 | 102 | 92 | 82 |
| MB 240-207720/8 | | 117 | 114 | 120 | 114 |
| LCS 240-207343/5 | | 97 | 94 | 91 | 80 |
| LCS 240-207343/7 | | 90 | 92 | 89 | 80 |
| LCS 240-207560/5 | | 98 | 98 | 92 | 81 |
| LCS 240-207560/7 | | 95 | 94 | 89 | 83 |
| LCS 240-207720/6 | | 105 | 100 | 104 | 96 |
| 240-57921-A-4-J MS | | 52 | 93 | 90 | 83 |
| 240-57972-C-3-C MS | | 101 | 94 | 101 | 98 |
| 240-57921-A-4-K MSD | | 49 | 93 | 90 | 84 |
| 240-57972-C-3-D MSD | | 98 | 95 | 101 | 100 |

| Surrogate | Acceptance Limits |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 37-132 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 58-123 |
| TOL = Toluene-d8 (Surr) | 67-125 |
| BFB = 4-Bromofluorobenzene (Surr) | 52-136 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Surrogate Recovery Report

**8260C Volatile Organic Compounds by GC/MS**

**Client Matrix: Water**

| Lab Sample ID | Client Sample ID | DBFM %Rec | DCA %Rec | TOL %Rec | BFB %Rec |
|---|---|---|---|---|---|
| 240-57930-12 | TB-111315 | 81 | 83 | 82 | 75 |
| MB 240-207096/6 | | 82 | 82 | 82 | 75 |
| LCS 240-207096/4 | | 90 | 90 | 90 | 85 |
| 200-30727-B-1 MS | | 81 | 81 | 84 | 80 |
| 200-30727-B-1 MSD | | 84 | 80 | 84 | 80 |

| Surrogate | Acceptance Limits |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 79-120 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 78-125 |
| TOL = Toluene-d8 (Surr) | 80-120 |
| BFB = 4-Bromofluorobenzene (Surr) | 61-120 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Surrogate Recovery Report

**8270D Semivolatile Organic Compounds (GC/MS)**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | 2FP %Rec | PHL %Rec | NBZ %Rec | FBP %Rec | TBP %Rec | TPH %Rec |
|---|---|---|---|---|---|---|---|
| 240-57930-1 | F-VS-UST-P-1 | 43 | 44 | 42 | 44 | 39 | 44 |
| 240-57930-2 | F-VS-UST-P-2 | 72 | 72 | 68 | 72 | 58 | 76 |
| 240-57930-3 | F-VS-UST-P-3 | 80 | 82 | 71 | 77 | 65 | 76 |
| 240-57930-4 | F-VS-UST-P-4 | 66 | 67 | 62 | 66 | 56 | 69 |
| 240-57930-5 | F-VB-UST-P-1 | 78 | 80 | 70 | 79 | 56 | 86 |
| 240-57930-6 | F-VB-UST-P-2 | 68 | 68 | 64 | 67 | 55 | 68 |
| 240-57930-7 | F-VB-UST-P-3 | 66 | 67 | 59 | 63 | 40 | 66 |
| 240-57930-8 | F-VB-UST-P-4 | 73 | 75 | 66 | 71 | 54 | 74 |
| 240-57930-9 | F-VB-UST-P-5 | 68 | 69 | 60 | 67 | 41 | 73 |
| 240-57930-10 | F-VB-UST-P-6 | 68 | 69 | 61 | 66 | 53 | 70 |
| 240-57930-11 | F-VS-UST-P-DUP | 68 | 68 | 65 | 66 | 62 | 66 |
| MB 240-206887/23-A | | 65 | 72 | 67 | 74 | 43 | 77 |
| MB 240-206908/21-A | | 70 | 73 | 67 | 70 | 30 | 74 |
| LCS 240-206887/24-A | | 80 | 80 | 73 | 77 | 69 | 78 |
| LCS 240-206908/22-A | | 70 | 71 | 70 | 70 | 42 | 71 |
| 240-57930-5 MS | F-VB-UST-P-1 MS | 78 | 78 | 73 | 74 | 72 | 76 |
| 240-57930-5 MSD | F-VB-UST-P-1 MSD | 73 | 74 | 71 | 72 | 64 | 75 |

| Surrogate | Acceptance Limits |
|---|---|
| 2FP = 2-Fluorophenol (Surr) | 24-110 |
| PHL = Phenol-d5 (Surr) | 26-110 |
| NBZ = Nitrobenzene-d5 (Surr) | 20-110 |
| FBP = 2-Fluorobiphenyl (Surr) | 24-110 |
| TBP = 2,4,6-Tribromophenol (Surr) | 10-110 |
| TPH = Terphenyl-d14 (Surr) | 36-110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Surrogate Recovery Report

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | TFT1 %Rec |
|---|---|---|
| 240-57930-1 | F-VS-UST-P-1 | 69 |
| 240-57930-2 | F-VS-UST-P-2 | 78 |
| 240-57930-3 | F-VS-UST-P-3 | 75 |
| 240-57930-4 | F-VS-UST-P-4 | 67 |
| 240-57930-5 | F-VB-UST-P-1 | 72 |
| 240-57930-6 | F-VB-UST-P-2 | 64 |
| 240-57930-7 | F-VB-UST-P-3 | 72 |
| 240-57930-8 | F-VB-UST-P-4 | 53 |
| 240-57930-9 | F-VB-UST-P-5 | 51 |
| 240-57930-10 | F-VB-UST-P-6 | 72 |
| 240-57930-11 | F-VS-UST-P-DUP | 70 |
| MB 240-206878/1-A | | 71 |
| MB 240-207078/1-A | | 79 |
| MB 240-207274/1-A | | 78 |
| LCS 240-206878/2-A | | 72 |
| LCS 240-207078/2-A | | 101 |
| LCS 240-207274/2-A | | 101 |
| 240-57930-11 MS | F-VS-UST-P-DUP MS | 87 |
| 240-57774-C-8-J MS | | 100 |
| 240-57930-11 MSD | F-VS-UST-P-DUP MSD | 73 |
| 240-57774-C-8-K MSD | | 100 |

| Surrogate | Acceptance Limits |
|---|---|
| TFT = Trifluorotoluene (Surr) | 40-139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Surrogate Recovery Report

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | OTPH1 %Rec |
|---|---|---|
| 240-57930-1 | F-VS-UST-P-1 | 82 |
| 240-57930-2 | F-VS-UST-P-2 | 95 |
| 240-57930-3 | F-VS-UST-P-3 | 75 |
| 240-57930-4 | F-VS-UST-P-4 | 82 |
| 240-57930-5 | F-VB-UST-P-1 | 76 |
| 240-57930-6 | F-VB-UST-P-2 | 79 |
| 240-57930-7 | F-VB-UST-P-3 | 69 |
| 240-57930-8 | F-VB-UST-P-4 | 90 |
| 240-57930-9 | F-VB-UST-P-5 | 81 |
| 240-57930-10 | F-VB-UST-P-6 | 83 |
| 240-57930-11 | F-VS-UST-P-DUP | 79 |
| MB 240-206884/23-A | | 77 |
| MB 240-207121/7-A | | 75 |
| LCS 240-206884/24-A | | 82 |
| LCS 240-207121/8-A | | 103 |
| 240-57930-11 MS | F-VS-UST-P-DUP MS | 93 |
| 240-57930-11 MSD | F-VS-UST-P-DUP MSD | 93 |

| Surrogate | Acceptance Limits |
|---|---|
| OTPH = o-Terphenyl (Surr) | 40-160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Surrogate Recovery Report

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | TCX2 %Rec | DCB2 %Rec |
|---|---|---|---|
| 240-57930-1 | F-VS-UST-P-1 | 83 | 72 |
| 240-57930-2 | F-VS-UST-P-2 | 61 | 70 |
| 240-57930-3 | F-VS-UST-P-3 | 67 | 90 |
| 240-57930-4 | F-VS-UST-P-4 | 78 | 141 |
| 240-57930-5 | F-VB-UST-P-1 | 83 | 82 |
| 240-57930-6 | F-VB-UST-P-2 | 70 | 66 |
| 240-57930-7 | F-VB-UST-P-3 | 67 | 66 |
| 240-57930-8 | F-VB-UST-P-4 | 83 | 79 |
| 240-57930-9 | F-VB-UST-P-5 | 80 | 71 |
| 240-57930-10 | F-VB-UST-P-6 | 72 | 71 |
| 240-57930-11 | F-VS-UST-P-DUP | 82 | 78 |
| MB 240-206917/21-A | | 74 | 79 |
| LCS 240-206917/22-A | | 75 | 77 |
| 240-57880-E-1-B MS | | 75 | 94 |
| 240-57880-E-1-C MSD | | 90 | 87 |

| Surrogate | Acceptance Limits |
|---|---|
| TCX = Tetrachloro-m-xylene | 13-134 |
| DCB = DCB Decachlorobiphenyl | 10-155 |

**Method Blank - Batch: 240-207096**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207096/6 | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3070.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1217 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1217 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.28 | U | 0.28 | 1.0 |
| 1,1,1-Trichloroethane | 0.44 | U | 0.44 | 1.0 |
| 1,1,2,2-Tetrachloroethane | 0.22 | U | 0.22 | 1.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloroethane | 0.30 | U | 0.30 | 1.0 |
| 1,1-Dichloroethene | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloropropene | 0.42 | U | 0.42 | 1.0 |
| 1,2,3-Trichlorobenzene | 0.37 | U | 0.37 | 1.0 |
| 1,2,3-Trichloropropane | 0.44 | U | 0.44 | 1.0 |
| 1,2,3-Trimethylbenzene | 0.47 | U | 0.47 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.32 | U | 0.32 | 1.0 |
| 1,2,4-Trimethylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromo-3-Chloropropane | 0.82 | U | 0.82 | 2.0 |
| 1,2-Dichlorobenzene | 0.25 | U | 0.25 | 1.0 |
| 1,2-Dichloroethane | 0.23 | U | 0.23 | 1.0 |
| 1,2-Dichloropropane | 0.25 | U | 0.25 | 1.0 |
| 1,3-Dichlorobenzene | 0.19 | U | 0.19 | 1.0 |
| 1,3-Dichloropropane | 0.19 | U | 0.19 | 1.0 |
| 1,4-Dichlorobenzene | 0.27 | U | 0.27 | 1.0 |
| 2,2-Dichloropropane | 0.26 | U | 0.26 | 1.0 |
| 2-Butanone (MEK) | 0.53 | U | 0.53 | 10 |
| 2-Chloroethyl vinyl ether | 0.59 | U | 0.59 | 10 |
| 2-Chlorotoluene | 0.40 | U | 0.40 | 1.0 |
| 2-Hexanone | 0.48 | U | 0.48 | 10 |
| 4-Chlorotoluene | 0.29 | U | 0.29 | 1.0 |
| p-Isopropyltoluene | 0.43 | U | 0.43 | 1.0 |
| 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 0.99 | 10 |
| Acetone | 0.94 | U | 0.94 | 10 |
| Benzene | 0.35 | U | 0.35 | 1.0 |
| Bromobenzene | 0.35 | U | 0.35 | 1.0 |
| Bromoform | 0.56 | U | 0.56 | 1.0 |
| Bromomethane | 0.44 | U | 0.44 | 1.0 |
| Carbon disulfide | 0.38 | U | 0.38 | 1.0 |
| Carbon tetrachloride | 0.43 | U | 0.43 | 1.0 |
| Chlorobenzene | 0.25 | U | 0.25 | 1.0 |
| Dibromochloromethane | 0.43 | U | 0.43 | 1.0 |
| Chloroethane | 0.32 | U | 0.32 | 1.0 |
| Chloroform | 0.25 | U | 0.25 | 1.0 |
| Chloromethane | 0.44 | U | 0.44 | 1.0 |
| cis-1,2-Dichloroethene | 0.26 | U | 0.26 | 1.0 |
| cis-1,3-Dichloropropene | 0.46 | U | 0.46 | 1.0 |
| Dibromomethane | 0.42 | U | 0.42 | 1.0 |
| Bromochloromethane | 0.50 | U | 0.50 | 1.0 |
| Dichlorodifluoromethane | 0.32 | U | 0.32 | 1.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 1.0 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-207096**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207096/6 | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3070.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1217 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1217 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| tert-Butyl alcohol (TBA) | 4.9 | U | 4.9 | 50 |
| Ethylbenzene | 0.25 | U | 0.25 | 1.0 |
| Hexachlorobutadiene | 0.35 | U | 0.35 | 1.0 |
| Diisopropyl ether | 0.50 | U | 0.50 | 10 |
| Isopropylbenzene | 0.35 | U | 0.35 | 1.0 |
| Methyl tert-butyl ether | 0.20 | U | 0.20 | 1.0 |
| Methylene Chloride | 1.18 | | 0.33 | 1.0 |
| m-Xylene & p-Xylene | 0.24 | U | 0.24 | 2.0 |
| Naphthalene | 0.45 | U | 0.45 | 1.0 |
| n-Butylbenzene | 0.31 | U | 0.31 | 1.0 |
| N-Propylbenzene | 0.40 | U | 0.40 | 1.0 |
| o-Xylene | 0.25 | U | 0.25 | 1.0 |
| sec-Butylbenzene | 0.48 | U | 0.48 | 1.0 |
| Styrene | 0.45 | U | 0.45 | 1.0 |
| Tert-amyl methyl ether | 0.30 | U | 0.30 | 5.0 |
| Ethyl t-butyl ether | 0.23 | U | 0.23 | 5.0 |
| Tetrachloroethene | 0.31 | U | 0.31 | 1.0 |
| Toluene | 0.23 | U | 0.23 | 1.0 |
| trans-1,2-Dichloroethene | 0.30 | U | 0.30 | 1.0 |
| trans-1,3-Dichloropropene | 0.56 | U | 0.56 | 1.0 |
| Trichloroethene | 0.22 | U | 0.22 | 1.0 |
| Trichlorofluoromethane | 0.49 | U | 0.49 | 1.0 |
| Vinyl acetate | 0.41 | U | 0.41 | 2.0 |
| Vinyl chloride | 0.29 | U | 0.29 | 1.0 |
| Xylenes, Total | 0.52 | U | 0.52 | 2.0 |
| tert-Butylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromoethane | 0.32 | U | 0.32 | 1.0 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 82 | 80 - 120 |
| Dibromofluoromethane (Surr) | 82 | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 75 | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 82 | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-207096**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207096/4 | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3068.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1134 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1134 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 10.0 | 9.46 | 95 | 80 - 120 | |
| 1,1,1-Trichloroethane | 10.0 | 9.11 | 91 | 77 - 123 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 10.0 | 100 | 71 - 123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 10.0 | 9.72 | 97 | 67 - 138 | |
| 1,1-Dichloroethane | 10.0 | 9.27 | 93 | 79 - 125 | |
| 1,1-Dichloroethene | 10.0 | 9.08 | 91 | 76 - 124 | |
| 1,1-Dichloropropene | 10.0 | 10.0 | 100 | 80 - 128 | |
| 1,2,3-Trichlorobenzene | 10.0 | 5.89 | 59 | 66 - 120 | * |
| 1,2,3-Trichloropropane | 10.0 | 10.0 | 100 | 72 - 126 | |
| 1,2,4-Trichlorobenzene | 10.0 | 6.37 | 64 | 61 - 120 | |
| 1,2,4-Trimethylbenzene | 10.0 | 9.26 | 93 | 76 - 120 | |
| 1,2-Dibromo-3-Chloropropane | 10.0 | 8.37 | 84 | 50 - 132 | |
| 1,2-Dichlorobenzene | 10.0 | 9.01 | 90 | 79 - 120 | |
| 1,2-Dichloroethane | 10.0 | 9.37 | 94 | 80 - 120 | |
| 1,2-Dichloropropane | 10.0 | 9.44 | 94 | 78 - 124 | |
| 1,3-Dichlorobenzene | 10.0 | 9.15 | 91 | 79 - 120 | |
| 1,3-Dichloropropane | 10.0 | 9.46 | 95 | 80 - 120 | |
| 1,4-Dichlorobenzene | 10.0 | 9.15 | 91 | 79 - 120 | |
| 2,2-Dichloropropane | 10.0 | 7.77 | 78 | 71 - 122 | |
| 2-Butanone (MEK) | 20.0 | 19.0 | 95 | 56 - 138 | |
| 2-Chloroethyl vinyl ether | 10.0 | 9.41 | 94 | 54 - 139 | J |
| 2-Chlorotoluene | 10.0 | 9.78 | 98 | 80 - 120 | |
| 2-Hexanone | 20.0 | 19.2 | 96 | 55 - 141 | |
| 4-Chlorotoluene | 10.0 | 9.88 | 99 | 80 - 120 | |
| p-Isopropyltoluene | 10.0 | 9.20 | 92 | 77 - 120 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 18.8 | 94 | 64 - 135 | |
| Acetone | 20.0 | 18.4 | 92 | 34 - 148 | |
| Benzene | 10.0 | 9.16 | 92 | 80 - 120 | |
| Bromobenzene | 10.0 | 9.75 | 98 | 80 - 120 | |
| Bromoform | 10.0 | 8.92 | 89 | 56 - 122 | |
| Bromomethane | 10.0 | 8.49 | 85 | 38 - 132 | |
| Carbon disulfide | 10.0 | 8.54 | 85 | 65 - 144 | |
| Carbon tetrachloride | 10.0 | 9.28 | 93 | 77 - 131 | |
| Chlorobenzene | 10.0 | 9.38 | 94 | 80 - 120 | |
| Dibromochloromethane | 10.0 | 9.64 | 96 | 74 - 120 | |
| Chloroethane | 10.0 | 8.47 | 85 | 36 - 126 | |
| Chloroform | 10.0 | 9.47 | 95 | 80 - 120 | |
| Chloromethane | 10.0 | 8.43 | 84 | 48 - 133 | |
| cis-1,2-Dichloroethene | 10.0 | 9.25 | 92 | 79 - 120 | |
| cis-1,3-Dichloropropene | 10.0 | 9.55 | 96 | 74 - 126 | |
| Dibromomethane | 10.0 | 9.34 | 93 | 80 - 120 | |
| Bromochloromethane | 10.0 | 9.25 | 93 | 80 - 120 | |
| Dichlorodifluoromethane | 10.0 | 6.79 | 68 | 23 - 136 | |
| Bromodichloromethane | 10.0 | 9.05 | 90 | 80 - 120 | |
| tert-Butyl alcohol (TBA) | 100 | 73.6 | 74 | 38 - 144 | |
| Ethylbenzene | 10.0 | 9.52 | 95 | 80 - 120 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-207096**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207096/4 | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3068.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1134 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1134 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Hexachlorobutadiene | 10.0 | 6.28 | 63 | 62 - 120 | |
| Isopropylbenzene | 10.0 | 9.49 | 95 | 77 - 120 | |
| Methyl tert-butyl ether | 10.0 | 8.83 | 88 | 69 - 121 | |
| Methylene Chloride | 10.0 | 10.9 | 109 | 77 - 129 | |
| m-Xylene & p-Xylene | 10.0 | 9.72 | 97 | 80 - 120 | |
| Naphthalene | 10.0 | 6.29 | 63 | 53 - 121 | |
| n-Butylbenzene | 10.0 | 8.80 | 88 | 74 - 120 | |
| N-Propylbenzene | 10.0 | 10.1 | 101 | 79 - 123 | |
| o-Xylene | 10.0 | 9.71 | 97 | 80 - 120 | |
| sec-Butylbenzene | 10.0 | 9.36 | 94 | 77 - 120 | |
| Styrene | 10.0 | 9.41 | 94 | 76 - 122 | |
| Tetrachloroethene | 10.0 | 9.47 | 95 | 78 - 121 | |
| Toluene | 10.0 | 9.21 | 92 | 80 - 120 | |
| trans-1,2-Dichloroethene | 10.0 | 9.18 | 92 | 80 - 124 | |
| trans-1,3-Dichloropropene | 10.0 | 9.95 | 100 | 75 - 131 | |
| Trichloroethene | 10.0 | 9.58 | 96 | 80 - 121 | |
| Trichlorofluoromethane | 10.0 | 9.63 | 96 | 61 - 133 | |
| Vinyl acetate | 10.0 | 7.50 | 75 | 27 - 160 | |
| Vinyl chloride | 10.0 | 8.72 | 87 | 52 - 121 | |
| Xylenes, Total | 20.0 | 19.4 | 97 | 80 - 120 | |
| tert-Butylbenzene | 10.0 | 9.70 | 97 | 77 - 121 | |
| 1,2-Dibromoethane | 10.0 | 9.91 | 99 | 80 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 90 | 80 - 120 |
| Dibromofluoromethane (Surr) | 90 | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 85 | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 90 | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207096**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 200-30727-B-1 MS | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3079.D |
| Dilution: | 50 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1535 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1535 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 200-30727-B-1 MSD | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3080.D |
| Dilution: | 50 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1557 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1557 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1,1,2-Tetrachloroethane | 97 | 93 | 72 - 120 | 4 | 22 | | |
| 1,1,1-Trichloroethane | 93 | 96 | 69 - 122 | 3 | 14 | | |
| 1,1,2,2-Tetrachloroethane | 97 | 97 | 61 - 130 | 0 | 18 | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 100 | 104 | 44 - 140 | 3 | 35 | | |
| 1,1-Dichloroethane | 95 | 96 | 73 - 124 | 0 | 14 | | |
| 1,1-Dichloroethene | 97 | 101 | 67 - 124 | 4 | 24 | | |
| 1,1-Dichloropropene | 105 | 103 | 81 - 126 | 2 | 11 | | |
| 1,2,3-Trichlorobenzene | 82 | 83 | 46 - 120 | 1 | 35 | | |
| 1,2,3-Trichloropropane | 100 | 98 | 58 - 135 | 1 | 25 | | |
| 1,2,4-Trichlorobenzene | 85 | 85 | 48 - 120 | 0 | 28 | | |
| 1,2,4-Trimethylbenzene | 95 | 94 | 64 - 122 | 2 | 14 | | |
| 1,2-Dibromo-3-Chloropropane | 76 | 77 | 42 - 130 | 2 | 24 | | |
| 1,2-Dichlorobenzene | 95 | 93 | 67 - 118 | 2 | 15 | | |
| 1,2-Dichloroethane | 98 | 97 | 74 - 125 | 0 | 24 | | |
| 1,2-Dichloropropane | 97 | 97 | 73 - 122 | 1 | 15 | | |
| 1,3-Dichlorobenzene | 94 | 94 | 65 - 120 | 0 | 15 | | |
| 1,3-Dichloropropane | 97 | 97 | 59 - 133 | 0 | 35 | | |
| 1,4-Dichlorobenzene | 94 | 93 | 66 - 120 | 1 | 16 | | |
| 2,2-Dichloropropane | 80 | 82 | 48 - 126 | 2 | 30 | | |
| 2-Butanone (MEK) | 95 | 100 | 49 - 132 | 4 | 19 | | |
| 2-Chloroethyl vinyl ether | 56 | 51 | 10 - 150 | 9 | 30 | J | J |
| 2-Chlorotoluene | 97 | 97 | 70 - 122 | 0 | 18 | | |
| 2-Hexanone | 95 | 98 | 49 - 142 | 3 | 27 | | |
| 4-Chlorotoluene | 100 | 97 | 74 - 120 | 3 | 15 | | |
| p-Isopropyltoluene | 95 | 95 | 61 - 123 | 0 | 27 | | |
| 4-Methyl-2-pentanone (MIBK) | 96 | 96 | 58 - 136 | 1 | 32 | | |
| Acetone | 100 | 107 | 32 - 126 | 7 | 28 | | |
| Benzene | 94 | 94 | 73 - 121 | 1 | 13 | | |
| Bromobenzene | 97 | 99 | 73 - 120 | 2 | 13 | | |
| Bromoform | 84 | 83 | 45 - 121 | 1 | 19 | | |
| Bromomethane | 80 | 84 | 26 - 136 | 4 | 35 | | |
| Carbon disulfide | 93 | 91 | 54 - 144 | 2 | 34 | | |
| Carbon tetrachloride | 95 | 97 | 65 - 129 | 2 | 20 | | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207096**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 200-30727-B-1 MS | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3079.D |
| Dilution: | 50 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1535 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1535 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 200-30727-B-1 MSD | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3080.D |
| Dilution: | 50 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1557 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1557 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Chlorobenzene | 96 | 95 | 72 - 120 | 1 | 15 | | |
| Dibromochloromethane | 92 | 93 | 62 - 122 | 1 | 19 | | |
| Chloroethane | 85 | 86 | 27 - 131 | 1 | 35 | | |
| Chloroform | 98 | 98 | 73 - 121 | 1 | 17 | | |
| Chloromethane | 71 | 76 | 39 - 134 | 6 | 20 | | |
| cis-1,2-Dichloroethene | 95 | 94 | 66 - 124 | 0 | 22 | | |
| cis-1,3-Dichloropropene | 94 | 97 | 60 - 120 | 3 | 21 | | |
| Dibromomethane | 96 | 96 | 63 - 134 | 0 | 35 | | |
| Bromochloromethane | 96 | 93 | 60 - 132 | 3 | 35 | | |
| Dichlorodifluoromethane | 72 | 71 | 14 - 137 | 2 | 34 | | |
| Bromodichloromethane | 92 | 92 | 72 - 120 | 1 | 19 | | |
| Ethylbenzene | 98 | 98 | 68 - 121 | 0 | 16 | | |
| Hexachlorobutadiene | 79 | 82 | 48 - 120 | 4 | 34 | | |
| Isopropylbenzene | 100 | 96 | 61 - 122 | 4 | 20 | | |
| Methyl tert-butyl ether | 89 | 91 | 61 - 121 | 3 | 12 | | |
| Methylene Chloride | 105 | 106 | 70 - 124 | 0 | 14 | | |
| m-Xylene & p-Xylene | 102 | 98 | 66 - 123 | 4 | 15 | | |
| Naphthalene | 100 | 86 | 38 - 128 | 14 | 35 | | |
| n-Butylbenzene | 95 | 93 | 54 - 126 | 2 | 33 | | |
| N-Propylbenzene | 103 | 102 | 63 - 129 | 1 | 29 | | |
| o-Xylene | 100 | 98 | 68 - 121 | 2 | 14 | | |
| sec-Butylbenzene | 98 | 96 | 61 - 123 | 2 | 29 | | |
| Styrene | 97 | 93 | 64 - 126 | 5 | 15 | | |
| Tetrachloroethene | 96 | 100 | 59 - 125 | 2 | 20 | | |
| Toluene | 97 | 96 | 72 - 122 | 1 | 15 | | |
| trans-1,2-Dichloroethene | 99 | 97 | 72 - 125 | 1 | 25 | | |
| trans-1,3-Dichloropropene | 95 | 99 | 58 - 132 | 4 | 22 | | |
| Trichloroethene | 103 | 103 | 61 - 129 | 0 | 14 | | |
| Trichlorofluoromethane | 91 | 96 | 49 - 133 | 5 | 25 | | |
| Vinyl acetate | 75 | 77 | 26 - 189 | 2 | 15 | | |
| Vinyl chloride | 80 | 84 | 44 - 122 | 3 | 35 | | |
| Xylenes, Total | 101 | 98 | 67 - 122 | 3 | 14 | | |
| tert-Butylbenzene | 97 | 97 | 63 - 122 | 0 | 24 | | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207096**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 200-30727-B-1 MS | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3079.D |
| Dilution: | 50 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1535 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1535 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 200-30727-B-1 MSD | Analysis Batch: | 240-207096 | Instrument ID: | A3UX10 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXX3080.D |
| Dilution: | 50 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 11/17/2015 1557 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1557 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,2-Dibromoethane | 99 | 97 | 69 - 125 | 2 | 24 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 84 | 84 | 80 - 120 |
| Dibromofluoromethane (Surr) | 81 | 84 | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 80 | 80 | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 81 | 80 | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207096**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 200-30727-B-1 MS | Units: ug/L | MSD Lab Sample ID: | 200-30727-B-1 MSD | |
| Client Matrix: | Water | | Client Matrix: | Water | |
| Dilution: | 50 | | Dilution: | 50 | |
| Analysis Date: | 11/17/2015 1535 | | Analysis Date: | 11/17/2015 1557 | |
| Prep Date: | 11/17/2015 1535 | | Prep Date: | 11/17/2015 1557 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 14 | U | 500 | 500 | 486 | | 467 | |
| 1,1,1-Trichloroethane | 22 | U | 500 | 500 | 466 | | 479 | |
| 1,1,2,2-Tetrachloroethane | 11 | U | 500 | 500 | 484 | | 485 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 23 | U | 500 | 500 | 502 | | 518 | |
| 1,1-Dichloroethane | 15 | U | 500 | 500 | 477 | | 479 | |
| 1,1-Dichloroethene | 23 | U | 500 | 500 | 485 | | 504 | |
| 1,1-Dichloropropene | 21 | U | 500 | 500 | 526 | | 514 | |
| 1,2,3-Trichlorobenzene | 19 | U | 500 | 500 | 409 | | 414 | |
| 1,2,3-Trichloropropane | 22 | U | 500 | 500 | 498 | | 491 | |
| 1,2,4-Trichlorobenzene | 16 | U | 500 | 500 | 426 | | 425 | |
| 1,2,4-Trimethylbenzene | 21 | U | 500 | 500 | 477 | | 470 | |
| 1,2-Dibromo-3-Chloropropane | 41 | U | 500 | 500 | 381 | | 387 | |
| 1,2-Dichlorobenzene | 13 | U | 500 | 500 | 475 | | 465 | |
| 1,2-Dichloroethane | 12 | U | 500 | 500 | 489 | | 487 | |
| 1,2-Dichloropropane | 13 | U | 500 | 500 | 483 | | 486 | |
| 1,3-Dichlorobenzene | 9.5 | U | 500 | 500 | 468 | | 469 | |
| 1,3-Dichloropropane | 9.5 | U | 500 | 500 | 483 | | 483 | |
| 1,4-Dichlorobenzene | 14 | U | 500 | 500 | 472 | | 465 | |
| 2,2-Dichloropropane | 13 | U | 500 | 500 | 402 | | 410 | |
| 2-Butanone (MEK) | 27 | U | 1000 | 1000 | 953 | | 997 | |
| 2-Chloroethyl vinyl ether | 30 | U | 500 | 500 | 278 | J | 254 | J |
| 2-Chlorotoluene | 20 | U | 500 | 500 | 484 | | 485 | |
| 2-Hexanone | 24 | U | 1000 | 1000 | 954 | | 985 | |
| 4-Chlorotoluene | 15 | U | 500 | 500 | 501 | | 486 | |
| p-Isopropyltoluene | 22 | U | 500 | 500 | 475 | | 477 | |
| 4-Methyl-2-pentanone (MIBK) | 50 | U | 1000 | 1000 | 957 | | 965 | |
| Acetone | 47 | U | 1000 | 1000 | 998 | | 1070 | |
| Benzene | 18 | U | 500 | 500 | 471 | | 469 | |
| Bromobenzene | 18 | U | 500 | 500 | 486 | | 494 | |
| Bromoform | 28 | U | 500 | 500 | 419 | | 416 | |
| Bromomethane | 22 | U | 500 | 500 | 402 | | 420 | |
| Carbon disulfide | 19 | U | 500 | 500 | 465 | | 455 | |
| Carbon tetrachloride | 22 | U | 500 | 500 | 475 | | 485 | |
| Chlorobenzene | 13 | U | 500 | 500 | 482 | | 477 | |
| Dibromochloromethane | 22 | U | 500 | 500 | 458 | | 464 | |
| Chloroethane | 16 | U | 500 | 500 | 426 | | 432 | |
| Chloroform | 13 | U | 500 | 500 | 488 | | 492 | |
| Chloromethane | 22 | U | 500 | 500 | 356 | | 379 | |
| cis-1,2-Dichloroethene | 1200 | | 500 | 500 | 1660 | | 1660 | |
| cis-1,3-Dichloropropene | 23 | U | 500 | 500 | 470 | | 483 | |
| Dibromomethane | 21 | U | 500 | 500 | 479 | | 478 | |
| Bromochloromethane | 25 | U | 500 | 500 | 481 | | 467 | |
| Dichlorodifluoromethane | 16 | U | 500 | 500 | 360 | | 354 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207096**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 200-30727-B-1 MS | Units: ug/L | MSD Lab Sample ID: | 200-30727-B-1 MSD | |
| Client Matrix: | Water | | Client Matrix: | Water | |
| Dilution: | 50 | | Dilution: | 50 | |
| Analysis Date: | 11/17/2015 1535 | | Analysis Date: | 11/17/2015 1557 | |
| Prep Date: | 11/17/2015 1535 | | Prep Date: | 11/17/2015 1557 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| Bromodichloromethane | 15 | U | 500 | 500 | 458 | 462 |
| Ethylbenzene | 13 | U | 500 | 500 | 489 | 489 |
| Hexachlorobutadiene | 18 | U | 500 | 500 | 395 | 410 |
| Isopropylbenzene | 18 | U | 500 | 500 | 498 | 479 |
| Methyl tert-butyl ether | 10 | U | 500 | 500 | 443 | 457 |
| Methylene Chloride | 52 | | 500 | 500 | 578 | 579 |
| m-Xylene & p-Xylene | 12 | U | 500 | 500 | 509 | 488 |
| Naphthalene | 23 | U | 500 | 500 | 498 | 431 |
| n-Butylbenzene | 16 | U | 500 | 500 | 473 | 463 |
| N-Propylbenzene | 20 | U | 500 | 500 | 516 | 510 |
| o-Xylene | 13 | U | 500 | 500 | 501 | 492 |
| sec-Butylbenzene | 24 | U | 500 | 500 | 489 | 480 |
| Styrene | 23 | U | 500 | 500 | 487 | 464 |
| Tetrachloroethene | 660 | | 500 | 500 | 1140 | 1160 |
| Toluene | 12 | U | 500 | 500 | 483 | 478 |
| trans-1,2-Dichloroethene | 15 | U | 500 | 500 | 493 | 486 |
| trans-1,3-Dichloropropene | 28 | U | 500 | 500 | 474 | 494 |
| Trichloroethene | 590 | | 500 | 500 | 1110 | 1110 |
| Trichlorofluoromethane | 25 | U | 500 | 500 | 455 | 479 |
| Vinyl acetate | 21 | U | 500 | 500 | 376 | 383 |
| Vinyl chloride | 290 | | 500 | 500 | 693 | 716 |
| Xylenes, Total | 26 | U | 1000 | 1000 | 1010 | 980 |
| tert-Butylbenzene | 21 | U | 500 | 500 | 484 | 485 |
| 1,2-Dibromoethane | 16 | U | 500 | 500 | 494 | 485 |

**Method Blank - Batch: 240-207343**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207343/6 | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146535.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/18/2015 1220 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.50 | U | 0.50 | 5.0 |
| 1,1,1-Trichloroethane | 0.65 | U | 0.65 | 5.0 |
| 1,1,2,2-Tetrachloroethane | 0.31 | U | 0.31 | 5.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 0.42 | 5.0 |
| 1,1-Dichloroethane | 0.34 | U | 0.34 | 5.0 |
| 1,1-Dichloroethene | 0.80 | U | 0.80 | 5.0 |
| 1,1-Dichloropropene | 0.37 | U | 0.37 | 5.0 |
| 1,2,3-Trichlorobenzene | 0.50 | U | 0.50 | 5.0 |
| 1,2,3-Trichloropropane | 0.78 | U | 0.78 | 5.0 |
| 1,2,3-Trimethylbenzene | 0.46 | U | 0.46 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.38 | U | 0.38 | 5.0 |
| 1,2,4-Trimethylbenzene | 0.58 | U | 0.58 | 5.0 |
| 1,2-Dibromo-3-Chloropropane | 2.7 | U | 2.7 | 10 |
| 1,2-Dichlorobenzene | 0.34 | U | 0.34 | 5.0 |
| 1,2-Dichloroethane | 0.46 | U | 0.46 | 5.0 |
| 1,2-Dichloropropane | 0.14 | U | 0.14 | 5.0 |
| 1,3-Dichlorobenzene | 0.69 | U | 0.69 | 5.0 |
| 1,3-Dichloropropane | 0.40 | U | 0.40 | 5.0 |
| 1,4-Dichlorobenzene | 0.28 | U | 0.28 | 5.0 |
| 2,2-Dichloropropane | 0.44 | U | 0.44 | 5.0 |
| 2-Butanone (MEK) | 1.1 | U | 1.1 | 20 |
| 2-Chloroethyl vinyl ether | 0.64 | U | 0.64 | 50 |
| 2-Chlorotoluene | 0.30 | U | 0.30 | 5.0 |
| 2-Hexanone | 0.68 | U | 0.68 | 20 |
| 4-Chlorotoluene | 0.26 | U | 0.26 | 5.0 |
| p-Isopropyltoluene | 0.41 | U | 0.41 | 5.0 |
| 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 1.2 | 20 |
| Acetone | 5.7 | U | 5.7 | 20 |
| Benzene | 0.90 | U | 0.90 | 5.0 |
| Bromobenzene | 0.33 | U | 0.33 | 5.0 |
| Bromoform | 0.24 | U | 0.24 | 5.0 |
| Bromomethane | 0.42 | U | 0.42 | 5.0 |
| Carbon disulfide | 0.59 | U | 0.59 | 5.0 |
| Carbon tetrachloride | 0.65 | U | 0.65 | 5.0 |
| Chlorobenzene | 0.53 | U | 0.53 | 5.0 |
| Dibromochloromethane | 0.37 | U | 0.37 | 5.0 |
| Chloroethane | 0.44 | U | 0.44 | 5.0 |
| Chloroform | 0.36 | U | 0.36 | 5.0 |
| Chloromethane | 0.79 | U | 0.79 | 5.0 |
| cis-1,2-Dichloroethene | 0.28 | U | 0.28 | 5.0 |
| cis-1,3-Dichloropropene | 0.71 | U | 0.71 | 5.0 |
| Dibromomethane | 0.42 | U | 0.42 | 5.0 |
| Bromochloromethane | 0.33 | U | 0.33 | 5.0 |
| Dichlorodifluoromethane | 0.28 | U | 0.28 | 5.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 5.0 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-207343**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207343/6 | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146535.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/18/2015 1220 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| tert-Butyl alcohol (TBA) | 5.7 | U | 5.7 | 200 |
| Ethylbenzene | 0.32 | U | 0.32 | 5.0 |
| Hexachlorobutadiene | 0.40 | U | 0.40 | 5.0 |
| Diisopropyl ether | 0.59 | U | 0.59 | 10 |
| Isopropylbenzene | 0.20 | U | 0.20 | 5.0 |
| Methyl tert-butyl ether | 0.43 | U | 0.43 | 5.0 |
| Methylene Chloride | 0.74 | U | 0.74 | 5.0 |
| m-Xylene & p-Xylene | 0.42 | U | 0.42 | 10 |
| Naphthalene | 0.35 | U | 0.35 | 5.0 |
| n-Butylbenzene | 0.59 | U | 0.59 | 5.0 |
| N-Propylbenzene | 0.60 | U | 0.60 | 5.0 |
| o-Xylene | 0.57 | U | 0.57 | 5.0 |
| sec-Butylbenzene | 0.91 | U | 0.91 | 5.0 |
| Styrene | 0.40 | U | 0.40 | 5.0 |
| Tert-amyl methyl ether | 0.43 | U | 0.43 | 5.0 |
| Ethyl t-butyl ether | 0.41 | U | 0.41 | 5.0 |
| Tetrachloroethene | 0.80 | U | 0.80 | 5.0 |
| Toluene | 0.298 | J | 0.27 | 5.0 |
| trans-1,2-Dichloroethene | 0.32 | U | 0.32 | 5.0 |
| trans-1,3-Dichloropropene | 0.29 | U | 0.29 | 5.0 |
| Trichloroethene | 0.38 | U | 0.38 | 5.0 |
| Trichlorofluoromethane | 0.29 | U | 0.29 | 5.0 |
| Vinyl acetate | 0.66 | U | 0.66 | 10 |
| Vinyl chloride | 0.30 | U | 0.30 | 5.0 |
| Xylenes, Total | 0.54 | U | 0.54 | 10 |
| tert-Butylbenzene | 0.67 | U | 0.67 | 5.0 |
| 1,2-Dibromoethane | 0.53 | U | 0.53 | 5.0 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 87 | 67 - 125 |
| Dibromofluoromethane (Surr) | 91 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 76 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 95 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-207343**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207343/5 | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146534.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/18/2015 1158 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 21.0 | 105 | 80 - 120 | |
| 1,1,1-Trichloroethane | 20.0 | 22.4 | 112 | 77 - 126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 17.6 | 88 | 77 - 123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 20.3 | 101 | 80 - 138 | |
| 1,1-Dichloroethane | 20.0 | 18.2 | 91 | 76 - 120 | |
| 1,1-Dichloroethene | 20.0 | 17.0 | 85 | 75 - 135 | |
| 1,1-Dichloropropene | 20.0 | 19.3 | 97 | 80 - 120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 18.7 | 93 | 61 - 121 | |
| 1,2,3-Trichloropropane | 20.0 | 20.0 | 100 | 80 - 126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 18.7 | 93 | 64 - 124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 18.7 | 94 | 80 - 129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 18.1 | 90 | 61 - 132 | |
| 1,2-Dichlorobenzene | 20.0 | 18.0 | 90 | 76 - 120 | |
| 1,2-Dichloroethane | 20.0 | 19.6 | 98 | 72 - 120 | |
| 1,2-Dichloropropane | 20.0 | 16.9 | 84 | 80 - 120 | |
| 1,3-Dichlorobenzene | 20.0 | 17.9 | 89 | 78 - 120 | |
| 1,3-Dichloropropane | 20.0 | 17.4 | 87 | 80 - 120 | |
| 1,4-Dichlorobenzene | 20.0 | 17.6 | 88 | 75 - 120 | |
| 2,2-Dichloropropane | 20.0 | 22.1 | 111 | 69 - 135 | |
| 2-Butanone (MEK) | 40.0 | 26.6 | 67 | 52 - 131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 12.5 | 62 | 28 - 173 | J |
| 2-Chlorotoluene | 20.0 | 19.3 | 97 | 78 - 120 | |
| 2-Hexanone | 40.0 | 30.6 | 76 | 64 - 136 | |
| 4-Chlorotoluene | 20.0 | 19.5 | 97 | 79 - 120 | |
| p-Isopropyltoluene | 20.0 | 19.3 | 97 | 77 - 131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 30.2 | 76 | 67 - 135 | |
| Acetone | 40.0 | 36.6 | 92 | 41 - 137 | |
| Benzene | 20.0 | 18.3 | 91 | 79 - 120 | |
| Bromobenzene | 20.0 | 19.7 | 99 | 80 - 120 | |
| Bromoform | 20.0 | 16.8 | 84 | 62 - 133 | |
| Bromomethane | 20.0 | 18.2 | 91 | 42 - 136 | |
| Carbon disulfide | 20.0 | 16.2 | 81 | 62 - 146 | |
| Carbon tetrachloride | 20.0 | 22.6 | 113 | 71 - 129 | |
| Chlorobenzene | 20.0 | 18.9 | 95 | 78 - 120 | |
| Dibromochloromethane | 20.0 | 18.3 | 91 | 72 - 127 | |
| Chloroethane | 20.0 | 17.0 | 85 | 58 - 120 | |
| Chloroform | 20.0 | 19.9 | 100 | 77 - 120 | |
| Chloromethane | 20.0 | 13.3 | 67 | 50 - 120 | |
| cis-1,2-Dichloroethene | 20.0 | 20.0 | 100 | 76 - 120 | |
| cis-1,3-Dichloropropene | 20.0 | 16.5 | 82 | 74 - 128 | |
| Dibromomethane | 20.0 | 19.1 | 95 | 80 - 120 | |
| Bromochloromethane | 20.0 | 20.6 | 103 | 79 - 120 | |
| Dichlorodifluoromethane | 20.0 | 16.4 | 82 | 26 - 120 | |
| Bromodichloromethane | 20.0 | 18.4 | 92 | 80 - 122 | |
| tert-Butyl alcohol (TBA) | 200 | 135 | 68 | 60 - 125 | J |
| Ethylbenzene | 20.0 | 19.3 | 97 | 79 - 120 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-207343**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207343/5 | Analysis Batch: | 240-207343 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146534.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/18/2015 1158 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Hexachlorobutadiene | 20.0 | 19.5 | 97 | 54 - 131 | |
| Isopropylbenzene | 20.0 | 19.6 | 98 | 76 - 122 | |
| Methyl tert-butyl ether | 20.0 | 18.0 | 90 | 49 - 165 | |
| Methylene Chloride | 20.0 | 15.3 | 76 | 75 - 120 | |
| m-Xylene & p-Xylene | 20.0 | 18.8 | 94 | 80 - 120 | |
| Naphthalene | 20.0 | 16.8 | 84 | 65 - 123 | |
| n-Butylbenzene | 20.0 | 18.9 | 95 | 68 - 135 | |
| N-Propylbenzene | 20.0 | 19.0 | 95 | 80 - 129 | |
| o-Xylene | 20.0 | 20.1 | 101 | 80 - 120 | |
| sec-Butylbenzene | 20.0 | 18.6 | 93 | 74 - 129 | |
| Styrene | 20.0 | 18.2 | 91 | 80 - 120 | |
| Tetrachloroethene | 20.0 | 20.1 | 101 | 79 - 120 | |
| Toluene | 20.0 | 19.5 | 97 | 75 - 120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.1 | 100 | 78 - 120 | |
| trans-1,3-Dichloropropene | 20.0 | 18.1 | 90 | 73 - 131 | |
| Trichloroethene | 20.0 | 20.3 | 102 | 79 - 120 | |
| Trichlorofluoromethane | 20.0 | 20.8 | 104 | 57 - 146 | |
| Vinyl acetate | 20.0 | 15.5 | 77 | 70 - 130 | |
| Vinyl chloride | 20.0 | 14.2 | 71 | 57 - 120 | |
| Xylenes, Total | 40.0 | 38.9 | 97 | 80 - 120 | |
| tert-Butylbenzene | 20.0 | 19.0 | 95 | 76 - 126 | |
| 1,2-Dibromoethane | 20.0 | 20.0 | 100 | 80 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 91 | 67 - 125 |
| Dibromofluoromethane (Surr) | 97 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 80 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 94 | 58 - 123 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 89 | 67 - 125 |
| Dibromofluoromethane (Surr) | 90 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 80 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 92 | 58 - 123 |

**Method Blank - Batch: 240-207560**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207560/6 | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146568.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/19/2015 1229 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.50 | U | 0.50 | 5.0 |
| 1,1,1-Trichloroethane | 0.65 | U | 0.65 | 5.0 |
| 1,1,2,2-Tetrachloroethane | 0.31 | U | 0.31 | 5.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 0.42 | 5.0 |
| 1,1-Dichloroethane | 0.34 | U | 0.34 | 5.0 |
| 1,1-Dichloroethene | 0.80 | U | 0.80 | 5.0 |
| 1,1-Dichloropropene | 0.37 | U | 0.37 | 5.0 |
| 1,2,3-Trichlorobenzene | 0.50 | U | 0.50 | 5.0 |
| 1,2,3-Trichloropropane | 0.78 | U | 0.78 | 5.0 |
| 1,2,3-Trimethylbenzene | 0.46 | U | 0.46 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.38 | U | 0.38 | 5.0 |
| 1,2,4-Trimethylbenzene | 0.58 | U | 0.58 | 5.0 |
| 1,2-Dibromo-3-Chloropropane | 2.7 | U | 2.7 | 10 |
| 1,2-Dichlorobenzene | 0.34 | U | 0.34 | 5.0 |
| 1,2-Dichloroethane | 0.46 | U | 0.46 | 5.0 |
| 1,2-Dichloropropane | 0.14 | U | 0.14 | 5.0 |
| 1,3-Dichlorobenzene | 0.69 | U | 0.69 | 5.0 |
| 1,3-Dichloropropane | 0.40 | U | 0.40 | 5.0 |
| 1,4-Dichlorobenzene | 0.28 | U | 0.28 | 5.0 |
| 2,2-Dichloropropane | 0.44 | U | 0.44 | 5.0 |
| 2-Butanone (MEK) | 1.1 | U | 1.1 | 20 |
| 2-Chloroethyl vinyl ether | 0.64 | U | 0.64 | 50 |
| 2-Chlorotoluene | 0.30 | U | 0.30 | 5.0 |
| 2-Hexanone | 0.68 | U | 0.68 | 20 |
| 4-Chlorotoluene | 0.26 | U | 0.26 | 5.0 |
| p-Isopropyltoluene | 0.41 | U | 0.41 | 5.0 |
| 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 1.2 | 20 |
| Acetone | 5.7 | U | 5.7 | 20 |
| Benzene | 0.90 | U | 0.90 | 5.0 |
| Bromobenzene | 0.33 | U | 0.33 | 5.0 |
| Bromoform | 0.24 | U | 0.24 | 5.0 |
| Bromomethane | 0.42 | U | 0.42 | 5.0 |
| Carbon disulfide | 0.59 | U | 0.59 | 5.0 |
| Carbon tetrachloride | 0.65 | U | 0.65 | 5.0 |
| Chlorobenzene | 0.53 | U | 0.53 | 5.0 |
| Dibromochloromethane | 0.37 | U | 0.37 | 5.0 |
| Chloroethane | 0.44 | U | 0.44 | 5.0 |
| Chloroform | 0.36 | U | 0.36 | 5.0 |
| Chloromethane | 0.79 | U | 0.79 | 5.0 |
| cis-1,2-Dichloroethene | 0.28 | U | 0.28 | 5.0 |
| cis-1,3-Dichloropropene | 0.71 | U | 0.71 | 5.0 |
| Dibromomethane | 0.42 | U | 0.42 | 5.0 |
| Bromochloromethane | 0.33 | U | 0.33 | 5.0 |
| Dichlorodifluoromethane | 0.28 | U | 0.28 | 5.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 5.0 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-207560**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207560/6 | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146568.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/19/2015 1229 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| tert-Butyl alcohol (TBA) | 5.7 | U | 5.7 | 200 |
| Ethylbenzene | 0.32 | U | 0.32 | 5.0 |
| Hexachlorobutadiene | 0.40 | U | 0.40 | 5.0 |
| Diisopropyl ether | 0.59 | U | 0.59 | 10 |
| Isopropylbenzene | 0.20 | U | 0.20 | 5.0 |
| Methyl tert-butyl ether | 0.43 | U | 0.43 | 5.0 |
| Methylene Chloride | 0.74 | U | 0.74 | 5.0 |
| m-Xylene & p-Xylene | 0.42 | U | 0.42 | 10 |
| Naphthalene | 0.35 | U | 0.35 | 5.0 |
| n-Butylbenzene | 0.59 | U | 0.59 | 5.0 |
| N-Propylbenzene | 0.60 | U | 0.60 | 5.0 |
| o-Xylene | 0.57 | U | 0.57 | 5.0 |
| sec-Butylbenzene | 0.91 | U | 0.91 | 5.0 |
| Styrene | 0.40 | U | 0.40 | 5.0 |
| Tert-amyl methyl ether | 0.43 | U | 0.43 | 5.0 |
| Ethyl t-butyl ether | 0.41 | U | 0.41 | 5.0 |
| Tetrachloroethene | 0.80 | U | 0.80 | 5.0 |
| Toluene | 0.348 | J | 0.27 | 5.0 |
| trans-1,2-Dichloroethene | 0.32 | U | 0.32 | 5.0 |
| trans-1,3-Dichloropropene | 0.29 | U | 0.29 | 5.0 |
| Trichloroethene | 0.38 | U | 0.38 | 5.0 |
| Trichlorofluoromethane | 0.29 | U | 0.29 | 5.0 |
| Vinyl acetate | 0.66 | U | 0.66 | 10 |
| Vinyl chloride | 0.30 | U | 0.30 | 5.0 |
| Xylenes, Total | 0.54 | U | 0.54 | 10 |
| tert-Butylbenzene | 0.67 | U | 0.67 | 5.0 |
| 1,2-Dibromoethane | 0.53 | U | 0.53 | 5.0 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 92 | 67 - 125 |
| Dibromofluoromethane (Surr) | 98 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 82 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 102 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-207560**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207560/5 | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146567.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/19/2015 1208 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 21.2 | 106 | 80 - 120 | |
| 1,1,1-Trichloroethane | 20.0 | 23.2 | 116 | 77 - 126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 18.7 | 93 | 77 - 123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 22.0 | 110 | 80 - 138 | |
| 1,1-Dichloroethane | 20.0 | 18.9 | 94 | 76 - 120 | |
| 1,1-Dichloroethene | 20.0 | 18.5 | 92 | 75 - 135 | |
| 1,1-Dichloropropene | 20.0 | 21.0 | 105 | 80 - 120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 19.1 | 96 | 61 - 121 | |
| 1,2,3-Trichloropropane | 20.0 | 18.4 | 92 | 80 - 126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 19.0 | 95 | 64 - 124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 19.3 | 97 | 80 - 129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 18.5 | 92 | 61 - 132 | |
| 1,2-Dichlorobenzene | 20.0 | 18.8 | 94 | 76 - 120 | |
| 1,2-Dichloroethane | 20.0 | 20.7 | 104 | 72 - 120 | |
| 1,2-Dichloropropane | 20.0 | 17.5 | 87 | 80 - 120 | |
| 1,3-Dichlorobenzene | 20.0 | 18.4 | 92 | 78 - 120 | |
| 1,3-Dichloropropane | 20.0 | 18.4 | 92 | 80 - 120 | |
| 1,4-Dichlorobenzene | 20.0 | 18.1 | 91 | 75 - 120 | |
| 2,2-Dichloropropane | 20.0 | 23.0 | 115 | 69 - 135 | |
| 2-Butanone (MEK) | 40.0 | 30.7 | 77 | 52 - 131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 13.8 | 69 | 28 - 173 | J |
| 2-Chlorotoluene | 20.0 | 19.4 | 97 | 78 - 120 | |
| 2-Hexanone | 40.0 | 32.4 | 81 | 64 - 136 | |
| 4-Chlorotoluene | 20.0 | 19.4 | 97 | 79 - 120 | |
| p-Isopropyltoluene | 20.0 | 19.9 | 99 | 77 - 131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 31.6 | 79 | 67 - 135 | |
| Acetone | 40.0 | 33.8 | 84 | 41 - 137 | |
| Benzene | 20.0 | 18.6 | 93 | 79 - 120 | |
| Bromobenzene | 20.0 | 20.2 | 101 | 80 - 120 | |
| Bromoform | 20.0 | 17.5 | 88 | 62 - 133 | |
| Bromomethane | 20.0 | 18.6 | 93 | 42 - 136 | |
| Carbon disulfide | 20.0 | 16.2 | 81 | 62 - 146 | |
| Carbon tetrachloride | 20.0 | 22.9 | 115 | 71 - 129 | |
| Chlorobenzene | 20.0 | 19.5 | 98 | 78 - 120 | |
| Dibromochloromethane | 20.0 | 18.4 | 92 | 72 - 127 | |
| Chloroethane | 20.0 | 18.5 | 92 | 58 - 120 | |
| Chloroform | 20.0 | 20.8 | 104 | 77 - 120 | |
| Chloromethane | 20.0 | 14.3 | 72 | 50 - 120 | |
| cis-1,2-Dichloroethene | 20.0 | 20.5 | 102 | 76 - 120 | |
| cis-1,3-Dichloropropene | 20.0 | 17.9 | 89 | 74 - 128 | |
| Dibromomethane | 20.0 | 21.3 | 107 | 80 - 120 | |
| Bromochloromethane | 20.0 | 20.9 | 105 | 79 - 120 | |
| Dichlorodifluoromethane | 20.0 | 15.4 | 77 | 26 - 120 | |
| Bromodichloromethane | 20.0 | 19.3 | 96 | 80 - 122 | |
| tert-Butyl alcohol (TBA) | 200 | 159 | 80 | 60 - 125 | J |
| Ethylbenzene | 20.0 | 20.0 | 100 | 79 - 120 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-207560**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207560/5 | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 146567.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/19/2015 1208 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Hexachlorobutadiene | 20.0 | 20.9 | 105 | 54 - 131 | |
| Isopropylbenzene | 20.0 | 20.5 | 102 | 76 - 122 | |
| Methyl tert-butyl ether | 20.0 | 18.8 | 94 | 49 - 165 | |
| Methylene Chloride | 20.0 | 15.3 | 76 | 75 - 120 | |
| m-Xylene & p-Xylene | 20.0 | 19.9 | 99 | 80 - 120 | |
| Naphthalene | 20.0 | 17.7 | 88 | 65 - 123 | |
| n-Butylbenzene | 20.0 | 19.4 | 97 | 68 - 135 | |
| N-Propylbenzene | 20.0 | 19.3 | 96 | 80 - 129 | |
| o-Xylene | 20.0 | 20.3 | 102 | 80 - 120 | |
| sec-Butylbenzene | 20.0 | 19.4 | 97 | 74 - 129 | |
| Styrene | 20.0 | 18.4 | 92 | 80 - 120 | |
| Tetrachloroethene | 20.0 | 20.7 | 104 | 79 - 120 | |
| Toluene | 20.0 | 19.9 | 100 | 75 - 120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.8 | 104 | 78 - 120 | |
| trans-1,3-Dichloropropene | 20.0 | 19.3 | 97 | 73 - 131 | |
| Trichloroethene | 20.0 | 20.4 | 102 | 79 - 120 | |
| Trichlorofluoromethane | 20.0 | 21.9 | 109 | 57 - 146 | |
| Vinyl acetate | 20.0 | 17.2 | 86 | 70 - 130 | |
| Vinyl chloride | 20.0 | 14.8 | 74 | 57 - 120 | |
| Xylenes, Total | 40.0 | 40.2 | 101 | 80 - 120 | |
| tert-Butylbenzene | 20.0 | 19.2 | 96 | 76 - 126 | |
| 1,2-Dibromoethane | 20.0 | 20.9 | 104 | 80 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 92 | 67 - 125 |
| Dibromofluoromethane (Surr) | 98 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 81 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 98 | 58 - 123 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 89 | 67 - 125 |
| Dibromofluoromethane (Surr) | 95 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 83 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 94 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207613**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57921-A-4-J MS | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-207613 | Lab File ID: | 146590.D |
| Dilution: | 1.0 | Leach Batch: | 240-206968 | Initial Weight/Volume: | 5.16 g |
| Analysis Date: | 11/19/2015 2022 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | 5 mL |
| Leach Date: | 11/16/2015 1230 | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57921-A-4-K MSD | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-207613 | Lab File ID: | 146591.D |
| Dilution: | 1.0 | Leach Batch: | 240-206968 | Initial Weight/Volume: | 5.24 g |
| Analysis Date: | 11/19/2015 2043 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | 5 mL |
| Leach Date: | 11/16/2015 1230 | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1,1-Trichloroethane | 99 | 101 | 51 - 128 | 1 | 30 | | |
| 1,1,2,2-Tetrachloroethane | 0 | 0 | 16 - 179 | NC | 30 | U F1 | U F1 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 104 | 87 | 50 - 147 | 19 | 30 | | |
| 1,1-Dichloroethane | 80 | 82 | 54 - 122 | 0 | 30 | | |
| 1,1-Dichloroethene | 122 | 114 | 49 - 157 | 8 | 30 | | |
| 1,2,4-Trichlorobenzene | 47 | 55 | 10 - 120 | 13 | 30 | | |
| 1,2-Dibromo-3-Chloropropane | 33 | 29 | 10 - 153 | 14 | 30 | J | J |
| 1,2-Dichlorobenzene | 57 | 66 | 17 - 122 | 13 | 30 | | |
| 1,2-Dichloroethane | 90 | 90 | 49 - 123 | 2 | 30 | | |
| 1,2-Dichloropropane | 75 | 79 | 61 - 120 | 4 | 30 | | |
| 1,3-Dichlorobenzene | 52 | 64 | 16 - 126 | 19 | 30 | | |
| 1,4-Dichlorobenzene | 51 | 59 | 15 - 121 | 13 | 30 | | |
| 2-Butanone (MEK) | 72 | 64 | 30 - 143 | 12 | 30 | | |
| 2-Hexanone | 71 | 67 | 37 - 147 | 6 | 30 | | |
| 4-Methyl-2-pentanone (MIBK) | 69 | 68 | 43 - 147 | 3 | 30 | | |
| Acetone | 110 | 94 | 24 - 140 | 9 | 30 | | |
| Benzene | 79 | 83 | 53 - 120 | 3 | 30 | | |
| Bromoform | 69 | 68 | 18 - 129 | 4 | 30 | | |
| Bromomethane | 61 | 59 | 33 - 130 | 6 | 30 | | |
| Carbon disulfide | 64 | 67 | 20 - 151 | 2 | 30 | | |
| Carbon tetrachloride | 96 | 100 | 32 - 137 | 2 | 30 | | |
| Chlorobenzene | 69 | 75 | 37 - 120 | 6 | 30 | | |
| Dibromochloromethane | 71 | 68 | 29 - 135 | 6 | 30 | | |
| Chloroethane | 78 | 84 | 45 - 120 | 5 | 30 | | |
| Chloroform | 90 | 93 | 53 - 120 | 2 | 30 | | |
| Chloromethane | 60 | 61 | 34 - 120 | 0 | 30 | | |
| cis-1,2-Dichloroethene | 85 | 88 | 50 - 120 | 2 | 30 | | |
| cis-1,3-Dichloropropene | 49 | 47 | 27 - 133 | 5 | 30 | | |
| Dichlorodifluoromethane | 68 | 69 | 17 - 120 | 1 | 30 | | |
| Bromodichloromethane | 70 | 68 | 35 - 132 | 4 | 30 | | |
| Ethylbenzene | 69 | 77 | 30 - 131 | 9 | 30 | | |
| Isopropylbenzene | 66 | 77 | 21 - 134 | 14 | 30 | | |
| Methyl tert-butyl ether | 86 | 86 | 51 - 157 | 2 | 30 | | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207613**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57921-A-4-J MS | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-207613 | Lab File ID: | 146590.D |
| Dilution: | 1.0 | Leach Batch: | 240-206968 | Initial Weight/Volume: | 5.16 g |
| Analysis Date: | 11/19/2015 2022 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | 5 mL |
| Leach Date: | 11/16/2015 1230 | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57921-A-4-K MSD | Analysis Batch: | 240-207560 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-207613 | Lab File ID: | 146591.D |
| Dilution: | 1.0 | Leach Batch: | 240-206968 | Initial Weight/Volume: | 5.24 g |
| Analysis Date: | 11/19/2015 2043 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/19/2015 1300 | | | | 5 mL |
| Leach Date: | 11/16/2015 1230 | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Methylene Chloride | 80 | 87 | 54 - 120 | 7 | 30 | | |
| m-Xylene & p-Xylene | 67 | 75 | 29 - 131 | 9 | 30 | | |
| o-Xylene | 71 | 79 | 29 - 134 | 9 | 30 | | |
| Styrene | 63 | 67 | 27 - 127 | 5 | 30 | | |
| Tetrachloroethene | 76 | 84 | 31 - 135 | 9 | 30 | | |
| Toluene | 76 | 82 | 39 - 129 | 5 | 30 | | |
| trans-1,2-Dichloroethene | 85 | 88 | 50 - 123 | 2 | 30 | | |
| trans-1,3-Dichloropropene | 61 | 57 | 28 - 137 | 8 | 30 | | |
| Trichloroethene | 156 | 163 | 10 - 177 | 3 | 30 | | |
| Trichlorofluoromethane | 93 | 98 | 36 - 142 | 4 | 30 | | |
| Vinyl chloride | 67 | 69 | 42 - 120 | 1 | 30 | | |
| Xylenes, Total | 69 | 77 | 30 - 131 | 9 | 30 | | |
| 1,2-Dibromoethane | 85 | 83 | 45 - 127 | 3 | 30 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 90 | 90 | 67 - 125 |
| Dibromofluoromethane (Surr) | 52 | 49 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 83 | 84 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 93 | 93 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207613**

**Method: 8260C**
**Preparation: 5030C**

| | | | |
|---|---|---|---|
| MS Lab Sample ID: | 240-57921-A-4-J MS | Units: ug/Kg | |
| Client Matrix: | Solid | | |
| Dilution: | 1.0 | | |
| Analysis Date: | 11/19/2015 2022 | | |
| Prep Date: | 11/19/2015 1300 | | |
| Leach Date: | 11/16/2015 1230 | | |

| | |
|---|---|
| MSD Lab Sample ID: | 240-57921-A-4-K MSD |
| Client Matrix: | Solid |
| Dilution: | 1.0 |
| Analysis Date: | 11/19/2015 2043 |
| Prep Date: | 11/19/2015 1300 |
| Leach Date: | 11/16/2015 1230 |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 0.65 | U | 20.2 | 19.9 | 20.0 | | 20.2 | |
| 1,1,2,2-Tetrachloroethane | 0.31 | U | 20.2 | 19.9 | 0.31 | U F1 | 0.31 | U F1 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 20.2 | 19.9 | 21.1 | | 17.4 | |
| 1,1-Dichloroethane | 0.34 | U | 20.2 | 19.9 | 16.3 | | 16.3 | |
| 1,1-Dichloroethene | 0.79 | U | 20.2 | 19.9 | 24.7 | | 22.7 | |
| 1,2,4-Trichlorobenzene | 0.38 | U | 20.2 | 19.9 | 9.56 | | 10.9 | |
| 1,2-Dibromo-3-Chloropropane | 2.7 | U | 20.2 | 19.9 | 6.61 | J | 5.77 | J |
| 1,2-Dichlorobenzene | 0.34 | U | 20.2 | 19.9 | 11.6 | | 13.2 | |
| 1,2-Dichloroethane | 0.46 | U | 20.2 | 19.9 | 18.3 | | 17.9 | |
| 1,2-Dichloropropane | 0.14 | U | 20.2 | 19.9 | 15.3 | | 15.8 | |
| 1,3-Dichlorobenzene | 0.69 | U | 20.2 | 19.9 | 10.5 | | 12.7 | |
| 1,4-Dichlorobenzene | 0.28 | U | 20.2 | 19.9 | 10.3 | | 11.8 | |
| 2-Butanone (MEK) | 2.9 | J | 40.5 | 39.9 | 32.2 | | 28.6 | |
| 2-Hexanone | 1.3 | J | 40.5 | 39.9 | 29.9 | | 28.2 | |
| 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 40.5 | 39.9 | 28.0 | | 27.1 | |
| Acetone | 34 | | 40.5 | 39.9 | 78.3 | | 71.4 | |
| Benzene | 0.89 | U | 20.2 | 19.9 | 16.0 | | 16.5 | |
| Bromoform | 0.24 | U | 20.2 | 19.9 | 14.0 | | 13.5 | |
| Bromomethane | 0.42 | U | 20.2 | 19.9 | 12.4 | | 11.7 | |
| Carbon disulfide | 2.2 | J | 20.2 | 19.9 | 15.1 | | 15.5 | |
| Carbon tetrachloride | 0.65 | U | 20.2 | 19.9 | 19.5 | | 20.0 | |
| Chlorobenzene | 0.53 | U | 20.2 | 19.9 | 14.1 | | 14.9 | |
| Dibromochloromethane | 0.37 | U | 20.2 | 19.9 | 14.4 | | 13.5 | |
| Chloroethane | 0.44 | U | 20.2 | 19.9 | 15.8 | | 16.7 | |
| Chloroform | 0.36 | U | 20.2 | 19.9 | 18.3 | | 18.6 | |
| Chloromethane | 0.78 | U | 20.2 | 19.9 | 12.2 | | 12.1 | |
| cis-1,2-Dichloroethene | 0.28 | U | 20.2 | 19.9 | 17.3 | | 17.6 | |
| cis-1,3-Dichloropropene | 0.71 | U | 20.2 | 19.9 | 9.86 | | 9.38 | |
| Dichlorodifluoromethane | 0.28 | U | 20.2 | 19.9 | 13.8 | | 13.7 | |
| Bromodichloromethane | 0.29 | U | 20.2 | 19.9 | 14.2 | | 13.6 | |
| Ethylbenzene | 0.32 | U | 20.2 | 19.9 | 14.0 | | 15.4 | |
| Isopropylbenzene | 0.20 | U | 20.2 | 19.9 | 13.3 | | 15.4 | |
| Methyl tert-butyl ether | 0.43 | U | 20.2 | 19.9 | 17.4 | | 17.2 | |
| Methylene Chloride | 0.73 | U | 20.2 | 19.9 | 16.2 | | 17.4 | |
| m-Xylene & p-Xylene | 0.42 | U | 20.2 | 19.9 | 13.6 | | 14.9 | |
| o-Xylene | 0.57 | U | 20.2 | 19.9 | 14.4 | | 15.7 | |
| Styrene | 0.40 | U | 20.2 | 19.9 | 12.7 | | 13.4 | |
| Tetrachloroethene | 0.79 | U | 20.2 | 19.9 | 15.3 | | 16.8 | |
| Toluene | 0.40 | J | 20.2 | 19.9 | 15.9 | | 16.7 | |
| trans-1,2-Dichloroethene | 0.32 | U | 20.2 | 19.9 | 17.3 | | 17.6 | |
| trans-1,3-Dichloropropene | 0.29 | U | 20.2 | 19.9 | 12.3 | | 11.4 | |
| Trichloroethene | 0.38 | U | 20.2 | 19.9 | 31.6 | | 32.5 | |
| Trichlorofluoromethane | 0.29 | U | 20.2 | 19.9 | 18.7 | | 19.5 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207613**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57921-A-4-J MS | Units: ug/Kg | MSD Lab Sample ID: | 240-57921-A-4-K MSD | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 11/19/2015 2022 | | Analysis Date: | 11/19/2015 2043 | |
| Prep Date: | 11/19/2015 1300 | | Prep Date: | 11/19/2015 1300 | |
| Leach Date: | 11/16/2015 1230 | | Leach Date: | 11/16/2015 1230 | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| Vinyl chloride | 0.30 | U | 20.2 | 19.9 | 13.6 | 13.7 |
| Xylenes, Total | 0.54 | U | 40.5 | 39.9 | 28.0 | 30.6 |
| 1,2-Dibromoethane | 0.53 | U | 20.2 | 19.9 | 17.1 | 16.6 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207709**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57972-C-3-C MS | Analysis Batch: | 240-207720 | Instrument ID: | A3UX8 |
| Client Matrix: | Solid | Prep Batch: | 240-207709 | Lab File ID: | UX80889.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.28 g |
| Analysis Date: | 11/20/2015 0841 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/20/2015 0452 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57972-C-3-D MSD | Analysis Batch: | 240-207720 | Instrument ID: | A3UX8 |
| Client Matrix: | Solid | Prep Batch: | 240-207709 | Lab File ID: | UX80890.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.04 g |
| Analysis Date: | 11/20/2015 0902 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/20/2015 0452 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | |
|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Acetone | 69 | 82 | 24 - 140 | 21 | 30 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 101 | 101 | 67 - 125 |
| Dibromofluoromethane (Surr) | 101 | 98 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 98 | 100 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 94 | 95 | 58 - 123 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207709**

**Method: 8260C**
**Preparation: 5030C**

| | | | |
|---|---|---|---|
| MS Lab Sample ID: | 240-57972-C-3-C MS | Units: ug/Kg | MSD Lab Sample ID: 240-57972-C-3-D MSD |
| Client Matrix: | Solid | | Client Matrix: Solid |
| Dilution: | 1.0 | | Dilution: 1.0 |
| Analysis Date: | 11/20/2015 0841 | | Analysis Date: 11/20/2015 0902 |
| Prep Date: | 11/20/2015 0452 | | Prep Date: 11/20/2015 0452 |
| Leach Date: | N/A | | Leach Date: N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| Acetone | 6.4 | U | 54.2 | 56.7 | 37.4 | 46.4 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-207720**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207720/8 | Analysis Batch: | 240-207720 | Instrument ID: | A3UX8 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | UX80879.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/20/2015 0508 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.50 | U | 0.50 | 5.0 |
| 1,1,1-Trichloroethane | 0.65 | U | 0.65 | 5.0 |
| 1,1,2,2-Tetrachloroethane | 0.31 | U | 0.31 | 5.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 0.42 | 5.0 |
| 1,1-Dichloroethane | 0.34 | U | 0.34 | 5.0 |
| 1,1-Dichloroethene | 0.80 | U | 0.80 | 5.0 |
| 1,1-Dichloropropene | 0.37 | U | 0.37 | 5.0 |
| 1,2,3-Trichlorobenzene | 0.50 | U | 0.50 | 5.0 |
| 1,2,3-Trichloropropane | 0.78 | U | 0.78 | 5.0 |
| 1,2,3-Trimethylbenzene | 0.46 | U | 0.46 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.38 | U | 0.38 | 5.0 |
| 1,2,4-Trimethylbenzene | 0.58 | U | 0.58 | 5.0 |
| 1,2-Dibromo-3-Chloropropane | 2.7 | U | 2.7 | 10 |
| 1,2-Dichlorobenzene | 0.34 | U | 0.34 | 5.0 |
| 1,2-Dichloroethane | 0.46 | U | 0.46 | 5.0 |
| 1,2-Dichloropropane | 0.14 | U | 0.14 | 5.0 |
| 1,3-Dichlorobenzene | 0.69 | U | 0.69 | 5.0 |
| 1,3-Dichloropropane | 0.40 | U | 0.40 | 5.0 |
| 1,4-Dichlorobenzene | 0.28 | U | 0.28 | 5.0 |
| 2,2-Dichloropropane | 0.44 | U | 0.44 | 5.0 |
| 2-Butanone (MEK) | 1.1 | U | 1.1 | 20 |
| 2-Chloroethyl vinyl ether | 0.64 | U | 0.64 | 50 |
| 2-Chlorotoluene | 0.30 | U | 0.30 | 5.0 |
| 2-Hexanone | 0.68 | U | 0.68 | 20 |
| 4-Chlorotoluene | 0.26 | U | 0.26 | 5.0 |
| p-Isopropyltoluene | 0.41 | U | 0.41 | 5.0 |
| 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 1.2 | 20 |
| Acetone | 6.42 | J | 5.7 | 20 |
| Benzene | 0.90 | U | 0.90 | 5.0 |
| Bromobenzene | 0.33 | U | 0.33 | 5.0 |
| Bromoform | 0.24 | U | 0.24 | 5.0 |
| Bromomethane | 0.42 | U | 0.42 | 5.0 |
| Carbon disulfide | 0.59 | U | 0.59 | 5.0 |
| Carbon tetrachloride | 0.65 | U | 0.65 | 5.0 |
| Chlorobenzene | 0.53 | U | 0.53 | 5.0 |
| Dibromochloromethane | 0.37 | U | 0.37 | 5.0 |
| Chloroethane | 0.44 | U | 0.44 | 5.0 |
| Chloroform | 0.36 | U | 0.36 | 5.0 |
| Chloromethane | 0.79 | U | 0.79 | 5.0 |
| cis-1,2-Dichloroethene | 0.28 | U | 0.28 | 5.0 |
| cis-1,3-Dichloropropene | 0.71 | U | 0.71 | 5.0 |
| Dibromomethane | 0.42 | U | 0.42 | 5.0 |
| Bromochloromethane | 0.33 | U | 0.33 | 5.0 |
| Dichlorodifluoromethane | 0.28 | U | 0.28 | 5.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 5.0 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-207720**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207720/8 | Analysis Batch: | 240-207720 | Instrument ID: | A3UX8 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | UX80879.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/20/2015 0508 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| tert-Butyl alcohol (TBA) | 5.7 | U | 5.7 | 200 |
| Ethylbenzene | 0.32 | U | 0.32 | 5.0 |
| Hexachlorobutadiene | 0.40 | U | 0.40 | 5.0 |
| Diisopropyl ether | 0.59 | U | 0.59 | 10 |
| Isopropylbenzene | 0.20 | U | 0.20 | 5.0 |
| Methyl tert-butyl ether | 0.43 | U | 0.43 | 5.0 |
| Methylene Chloride | 0.74 | U | 0.74 | 5.0 |
| m-Xylene & p-Xylene | 0.42 | U | 0.42 | 10 |
| Naphthalene | 0.35 | U | 0.35 | 5.0 |
| n-Butylbenzene | 0.59 | U | 0.59 | 5.0 |
| N-Propylbenzene | 0.60 | U | 0.60 | 5.0 |
| o-Xylene | 0.57 | U | 0.57 | 5.0 |
| sec-Butylbenzene | 0.91 | U | 0.91 | 5.0 |
| Styrene | 0.40 | U | 0.40 | 5.0 |
| Tert-amyl methyl ether | 0.43 | U | 0.43 | 5.0 |
| Ethyl t-butyl ether | 0.41 | U | 0.41 | 5.0 |
| Tetrachloroethene | 0.80 | U | 0.80 | 5.0 |
| Toluene | 0.27 | U | 0.27 | 5.0 |
| trans-1,2-Dichloroethene | 0.32 | U | 0.32 | 5.0 |
| trans-1,3-Dichloropropene | 0.29 | U | 0.29 | 5.0 |
| Trichloroethene | 0.38 | U | 0.38 | 5.0 |
| Trichlorofluoromethane | 0.29 | U | 0.29 | 5.0 |
| Vinyl acetate | 0.66 | U | 0.66 | 10 |
| Vinyl chloride | 0.30 | U | 0.30 | 5.0 |
| Xylenes, Total | 0.54 | U | 0.54 | 10 |
| tert-Butylbenzene | 0.67 | U | 0.67 | 5.0 |
| 1,2-Dibromoethane | 0.53 | U | 0.53 | 5.0 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 120 | 67 - 125 |
| Dibromofluoromethane (Surr) | 117 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 114 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 114 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-207720**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207720/6 | Analysis Batch: | 240-207720 | Instrument ID: | A3UX8 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | UX80877.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/20/2015 0426 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 50.0 | 49.7 | 99 | 80 - 120 | |
| 1,1,1-Trichloroethane | 50.0 | 49.6 | 99 | 77 - 126 | |
| 1,1,2,2-Tetrachloroethane | 50.0 | 49.8 | 100 | 77 - 123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 50.0 | 51.8 | 104 | 80 - 138 | |
| 1,1-Dichloroethane | 50.0 | 48.9 | 98 | 76 - 120 | |
| 1,1-Dichloroethene | 50.0 | 48.3 | 97 | 75 - 135 | |
| 1,1-Dichloropropene | 50.0 | 52.2 | 104 | 80 - 120 | |
| 1,2,3-Trichlorobenzene | 50.0 | 45.0 | 90 | 61 - 121 | |
| 1,2,3-Trichloropropane | 50.0 | 52.8 | 106 | 80 - 126 | |
| 1,2,4-Trichlorobenzene | 50.0 | 46.3 | 93 | 64 - 124 | |
| 1,2,4-Trimethylbenzene | 50.0 | 49.4 | 99 | 80 - 129 | |
| 1,2-Dibromo-3-Chloropropane | 50.0 | 48.6 | 97 | 61 - 132 | |
| 1,2-Dichlorobenzene | 50.0 | 46.6 | 93 | 76 - 120 | |
| 1,2-Dichloroethane | 50.0 | 49.9 | 100 | 72 - 120 | |
| 1,2-Dichloropropane | 50.0 | 47.1 | 94 | 80 - 120 | |
| 1,3-Dichlorobenzene | 50.0 | 45.9 | 92 | 78 - 120 | |
| 1,3-Dichloropropane | 50.0 | 46.0 | 92 | 80 - 120 | |
| 1,4-Dichlorobenzene | 50.0 | 45.5 | 91 | 75 - 120 | |
| 2,2-Dichloropropane | 50.0 | 44.2 | 88 | 69 - 135 | |
| 2-Butanone (MEK) | 100 | 98.6 | 99 | 52 - 131 | |
| 2-Chloroethyl vinyl ether | 50.0 | NC | NaN | 28 - 173 | |
| 2-Chlorotoluene | 50.0 | 48.0 | 96 | 78 - 120 | |
| 2-Hexanone | 100 | 102 | 102 | 64 - 136 | |
| 4-Chlorotoluene | 50.0 | 45.9 | 92 | 79 - 120 | |
| p-Isopropyltoluene | 50.0 | 50.3 | 101 | 77 - 131 | |
| 4-Methyl-2-pentanone (MIBK) | 100 | 107 | 107 | 67 - 135 | |
| Acetone | 100 | 111 | 111 | 41 - 137 | |
| Benzene | 50.0 | 46.8 | 94 | 79 - 120 | |
| Bromobenzene | 50.0 | 46.6 | 93 | 80 - 120 | |
| Bromoform | 50.0 | 49.8 | 100 | 62 - 133 | |
| Bromomethane | 50.0 | 56.1 | 112 | 42 - 136 | |
| Carbon disulfide | 50.0 | 50.4 | 101 | 62 - 146 | |
| Carbon tetrachloride | 50.0 | 53.3 | 107 | 71 - 129 | |
| Chlorobenzene | 50.0 | 45.5 | 91 | 78 - 120 | |
| Dibromochloromethane | 50.0 | 52.7 | 105 | 72 - 127 | |
| Chloroethane | 50.0 | 49.4 | 99 | 58 - 120 | |
| Chloroform | 50.0 | 49.5 | 99 | 77 - 120 | |
| Chloromethane | 50.0 | 50.8 | 102 | 50 - 120 | |
| cis-1,2-Dichloroethene | 50.0 | 48.3 | 97 | 76 - 120 | |
| cis-1,3-Dichloropropene | 50.0 | 48.3 | 97 | 74 - 128 | |
| Dibromomethane | 50.0 | 49.9 | 100 | 80 - 120 | |
| Bromochloromethane | 50.0 | 49.5 | 99 | 79 - 120 | |
| Dichlorodifluoromethane | 50.0 | 52.4 | 105 | 26 - 120 | |
| Bromodichloromethane | 50.0 | 50.4 | 101 | 80 - 122 | |
| tert-Butyl alcohol (TBA) | 500 | 440 | 88 | 60 - 125 | |
| Ethylbenzene | 50.0 | 45.5 | 91 | 79 - 120 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-207720**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207720/6 | Analysis Batch: | 240-207720 | Instrument ID: | A3UX8 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | UX80877.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/20/2015 0426 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Hexachlorobutadiene | 50.0 | 48.6 | 97 | 54 - 131 | |
| Isopropylbenzene | 50.0 | 50.1 | 100 | 76 - 122 | |
| Methyl tert-butyl ether | 50.0 | 47.8 | 96 | 49 - 165 | |
| Methylene Chloride | 50.0 | 52.3 | 105 | 75 - 120 | |
| m-Xylene & p-Xylene | 50.0 | 46.2 | 92 | 80 - 120 | |
| Naphthalene | 50.0 | 52.0 | 104 | 65 - 123 | |
| n-Butylbenzene | 50.0 | 50.0 | 100 | 68 - 135 | |
| N-Propylbenzene | 50.0 | 48.6 | 97 | 80 - 129 | |
| o-Xylene | 50.0 | 47.8 | 96 | 80 - 120 | |
| sec-Butylbenzene | 50.0 | 51.0 | 102 | 74 - 129 | |
| Styrene | 50.0 | 45.9 | 92 | 80 - 120 | |
| Tetrachloroethene | 50.0 | 45.4 | 91 | 79 - 120 | |
| Toluene | 50.0 | 45.5 | 91 | 75 - 120 | |
| trans-1,2-Dichloroethene | 50.0 | 48.7 | 97 | 78 - 120 | |
| trans-1,3-Dichloropropene | 50.0 | 42.9 | 86 | 73 - 131 | |
| Trichloroethene | 50.0 | 48.9 | 98 | 79 - 120 | |
| Trichlorofluoromethane | 50.0 | 50.8 | 102 | 57 - 146 | |
| Vinyl acetate | 50.0 | 39.2 | 78 | 70 - 130 | |
| Vinyl chloride | 50.0 | 49.9 | 100 | 57 - 120 | |
| Xylenes, Total | 100 | 94.0 | 94 | 80 - 120 | |
| tert-Butylbenzene | 50.0 | 55.3 | 111 | 76 - 126 | |
| 1,2-Dibromoethane | 50.0 | 48.6 | 97 | 80 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 104 | 67 - 125 |
| Dibromofluoromethane (Surr) | 105 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 96 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 100 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-206887**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206887/23-A | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206887 | Lab File ID: | 51119007.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/19/2015 1150 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1'-Biphenyl | 3.5 | U | 3.5 | 50 |
| bis (2-chloroisopropyl) ether | 9.5 | U | 9.5 | 100 |
| 2,4,5-Trichlorophenol | 25 | U | 25 | 150 |
| 2,4,6-Trichlorophenol | 8.9 | U | 8.9 | 150 |
| 2,4-Dichlorophenol | 20 | U | 20 | 150 |
| 2,4-Dimethylphenol | 20 | U | 20 | 150 |
| 2,4-Dinitrophenol | 21 | U | 21 | 330 |
| 2,4-Dinitrotoluene | 17 | U | 17 | 200 |
| 2,6-Dinitrotoluene | 21 | U | 21 | 200 |
| 2-Chloronaphthalene | 0.45 | U | 0.45 | 50 |
| 2-Chlorophenol | 8.2 | U | 8.2 | 50 |
| 2-Methylnaphthalene | 0.50 | U | 0.50 | 6.7 |
| 2-Methylphenol | 11 | U | 11 | 200 |
| 2-Nitroaniline | 9.1 | U | 9.1 | 200 |
| 2-Nitrophenol | 8.3 | U | 8.3 | 50 |
| 3,3'-Dichlorobenzidine | 18 | U | 18 | 100 |
| 3-Nitroaniline | 16 | U | 16 | 200 |
| 4,6-Dinitro-2-methylphenol | 9.2 | U | 9.2 | 150 |
| 4-Bromophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Chloro-3-methylphenol | 21 | U | 21 | 150 |
| 4-Chloroaniline | 17 | U | 17 | 150 |
| 4-Chlorophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Nitroaniline | 26 | U | 26 | 200 |
| 4-Nitrophenol | 17 | U | 17 | 330 |
| Acenaphthene | 0.76 | U | 0.76 | 6.7 |
| Acenaphthylene | 0.35 | U | 0.35 | 6.7 |
| Acetophenone | 9.2 | U | 9.2 | 100 |
| Anthracene | 0.78 | U | 0.78 | 6.7 |
| Atrazine | 9.1 | U | 9.1 | 200 |
| Benzaldehyde | 12 | U | 12 | 100 |
| Benzo[a]anthracene | 0.63 | U | 0.63 | 6.7 |
| Benzo[a]pyrene | 0.64 | U | 0.64 | 6.7 |
| Benzo[b]fluoranthene | 0.59 | U | 0.59 | 6.7 |
| Benzo[g,h,i]perylene | 0.35 | U | 0.35 | 6.7 |
| Benzo[k]fluoranthene | 0.68 | U | 0.68 | 6.7 |
| Bis(2-chloroethoxy)methane | 22 | U | 22 | 100 |
| Bis(2-chloroethyl)ether | 2.0 | U | 2.0 | 100 |
| Bis(2-ethylhexyl) phthalate | 34.7 | J | 19 | 70 |
| Butyl benzyl phthalate | 10 | U | 10 | 70 |
| Caprolactam | 37 | U | 37 | 330 |
| Carbazole | 27 | U | 27 | 50 |
| Chrysene | 1.1 | U | 1.1 | 6.7 |
| Dibenz(a,h)anthracene | 0.66 | U | 0.66 | 6.7 |
| Dibenzofuran | 0.66 | U | 0.66 | 50 |
| Diethyl phthalate | 16 | U | 16 | 70 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-206887**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206887/23-A | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206887 | Lab File ID: | 51119007.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/19/2015 1150 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Dimethyl phthalate | 17 | U | 17 | 70 |
| Di-n-butyl phthalate | 15 | U | 15 | 70 |
| Di-n-octyl phthalate | 7.9 | U | 7.9 | 70 |
| Fluoranthene | 0.55 | U | 0.55 | 6.7 |
| Fluorene | 0.53 | U | 0.53 | 6.7 |
| Hexachlorobenzene | 2.1 | U | 2.1 | 6.7 |
| Hexachlorobutadiene | 5.6 | U | 5.6 | 50 |
| Hexachlorocyclopentadiene | 8.1 | U | 8.1 | 330 |
| Hexachloroethane | 9.0 | U | 9.0 | 50 |
| Indeno[1,2,3-cd]pyrene | 0.35 | U | 0.35 | 6.7 |
| Isophorone | 13 | U | 13 | 50 |
| N-Nitrosodi-n-propylamine | 6.3 | U | 6.3 | 50 |
| N-Nitrosodiphenylamine | 21 | U | 21 | 50 |
| Naphthalene | 0.82 | U | 0.82 | 6.7 |
| Nitrobenzene | 2.2 | U | 2.2 | 100 |
| Pentachlorophenol | 9.1 | U | 9.1 | 150 |
| Phenanthrene | 0.73 | U | 0.73 | 6.7 |
| Phenol | 7.3 | U | 7.3 | 50 |
| Pyrene | 0.44 | U | 0.44 | 6.7 |
| 3 & 4 Methylphenol | 20 | U | 20 | 400 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 77 | 36 - 110 |
| Phenol-d5 (Surr) | 72 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 67 | 20 - 110 |
| 2-Fluorophenol (Surr) | 65 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 74 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 43 | 10 - 110 |

**Lab Control Sample - Batch: 240-206887**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206887/24-A | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206887 | Lab File ID: | 51119008.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/19/2015 1213 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 486 | 73 | 35 - 110 | |
| bis (2-chloroisopropyl) ether | 667 | 490 | 73 | 29 - 110 | |
| 2,4,5-Trichlorophenol | 667 | 484 | 73 | 25 - 110 | |
| 2,4,6-Trichlorophenol | 667 | 468 | 70 | 12 - 110 | |
| 2,4-Dichlorophenol | 667 | 499 | 75 | 39 - 110 | |
| 2,4-Dimethylphenol | 667 | 428 | 64 | 29 - 110 | |
| 2,4-Dinitrophenol | 1330 | 583 | 44 | 10 - 110 | |
| 2,4-Dinitrotoluene | 667 | 537 | 81 | 48 - 110 | |
| 2,6-Dinitrotoluene | 667 | 539 | 81 | 45 - 110 | |
| 2-Chloronaphthalene | 667 | 491 | 74 | 32 - 110 | |
| 2-Chlorophenol | 667 | 509 | 76 | 37 - 110 | |
| 2-Methylnaphthalene | 667 | 503 | 75 | 36 - 110 | |
| 2-Methylphenol | 667 | 500 | 75 | 41 - 110 | |
| 2-Nitroaniline | 667 | 507 | 76 | 45 - 110 | |
| 2-Nitrophenol | 667 | 509 | 76 | 34 - 110 | |
| 3,3'-Dichlorobenzidine | 1330 | 770 | 58 | 28 - 110 | |
| 3-Nitroaniline | 667 | 450 | 68 | 44 - 110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 727 | 55 | 10 - 110 | |
| 4-Bromophenyl phenyl ether | 667 | 539 | 81 | 39 - 110 | |
| 4-Chloro-3-methylphenol | 667 | 518 | 78 | 48 - 110 | |
| 4-Chloroaniline | 667 | 358 | 54 | 30 - 110 | |
| 4-Chlorophenyl phenyl ether | 667 | 513 | 77 | 40 - 110 | |
| 4-Nitroaniline | 667 | 534 | 80 | 48 - 110 | |
| 4-Nitrophenol | 1330 | 929 | 70 | 28 - 110 | |
| Acenaphthene | 667 | 487 | 73 | 38 - 110 | |
| Acenaphthylene | 667 | 498 | 75 | 40 - 110 | |
| Acetophenone | 667 | 491 | 74 | 40 - 110 | |
| Anthracene | 667 | 510 | 77 | 48 - 110 | |
| Atrazine | 1330 | 1060 | 79 | 54 - 120 | |
| Benzaldehyde | 1330 | 1010 | 76 | 32 - 110 | |
| Benzo[a]anthracene | 667 | 517 | 78 | 50 - 110 | |
| Benzo[a]pyrene | 667 | 509 | 76 | 44 - 110 | |
| Benzo[b]fluoranthene | 667 | 509 | 76 | 43 - 110 | |
| Benzo[g,h,i]perylene | 667 | 518 | 78 | 51 - 110 | |
| Benzo[k]fluoranthene | 667 | 530 | 79 | 38 - 105 | |
| Bis(2-chloroethoxy)methane | 667 | 526 | 79 | 32 - 110 | |
| Bis(2-chloroethyl)ether | 667 | 515 | 77 | 34 - 110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 607 | 91 | 50 - 110 | |
| Butyl benzyl phthalate | 667 | 545 | 82 | 51 - 110 | |
| Caprolactam | 1330 | 1100 | 82 | 44 - 114 | |
| Carbazole | 667 | 530 | 80 | 50 - 110 | |
| Chrysene | 667 | 538 | 81 | 50 - 110 | |
| Dibenz(a,h)anthracene | 667 | 504 | 76 | 51 - 110 | |
| Dibenzofuran | 667 | 499 | 75 | 43 - 110 | |
| Diethyl phthalate | 667 | 544 | 82 | 52 - 110 | |
| Dimethyl phthalate | 667 | 523 | 79 | 50 - 110 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-206887**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206887/24-A | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206887 | Lab File ID: | 51119008.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/19/2015 1213 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0848 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | 667 | 549 | 82 | 51 - 110 | |
| Di-n-octyl phthalate | 667 | 515 | 77 | 48 - 110 | |
| Fluoranthene | 667 | 528 | 79 | 51 - 110 | |
| Fluorene | 667 | 491 | 74 | 46 - 110 | |
| Hexachlorobenzene | 667 | 518 | 78 | 43 - 110 | |
| Hexachlorobutadiene | 667 | 490 | 73 | 29 - 110 | |
| Hexachlorocyclopentadiene | 667 | 320 | 48 | 12 - 110 | J |
| Hexachloroethane | 667 | 473 | 71 | 30 - 110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 515 | 77 | 50 - 110 | |
| Isophorone | 667 | 527 | 79 | 36 - 110 | |
| N-Nitrosodi-n-propylamine | 667 | 513 | 77 | 38 - 110 | |
| N-Nitrosodiphenylamine | 1330 | 1030 | 77 | 46 - 110 | |
| Naphthalene | 667 | 505 | 76 | 36 - 110 | |
| Nitrobenzene | 667 | 508 | 76 | 32 - 110 | |
| Pentachlorophenol | 1330 | 665 | 50 | 10 - 110 | |
| Phenanthrene | 667 | 516 | 77 | 49 - 110 | |
| Phenol | 667 | 509 | 76 | 38 - 110 | |
| Pyrene | 667 | 507 | 76 | 49 - 110 | |
| 3 & 4 Methylphenol | 667 | 501 | 75 | 40 - 110 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 78 | 36 - 110 |
| Phenol-d5 (Surr) | 80 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 73 | 20 - 110 |
| 2-Fluorophenol (Surr) | 80 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 77 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 69 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-206908**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206908/21-A | Analysis Batch: | 240-207287 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206908 | Lab File ID: | 51118003.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/18/2015 0851 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1'-Biphenyl | 3.5 | U | 3.5 | 50 |
| bis (2-chloroisopropyl) ether | 9.5 | U | 9.5 | 100 |
| 2,4,5-Trichlorophenol | 25 | U | 25 | 150 |
| 2,4,6-Trichlorophenol | 8.9 | U | 8.9 | 150 |
| 2,4-Dichlorophenol | 20 | U | 20 | 150 |
| 2,4-Dimethylphenol | 20 | U | 20 | 150 |
| 2,4-Dinitrophenol | 21 | U | 21 | 330 |
| 2,4-Dinitrotoluene | 17 | U | 17 | 200 |
| 2,6-Dinitrotoluene | 21 | U | 21 | 200 |
| 2-Chloronaphthalene | 0.45 | U | 0.45 | 50 |
| 2-Chlorophenol | 8.2 | U | 8.2 | 50 |
| 2-Methylnaphthalene | 0.50 | U | 0.50 | 6.7 |
| 2-Methylphenol | 11 | U | 11 | 200 |
| 2-Nitroaniline | 9.1 | U | 9.1 | 200 |
| 2-Nitrophenol | 8.3 | U | 8.3 | 50 |
| 3,3'-Dichlorobenzidine | 18 | U | 18 | 100 |
| 3-Nitroaniline | 16 | U | 16 | 200 |
| 4,6-Dinitro-2-methylphenol | 9.2 | U | 9.2 | 150 |
| 4-Bromophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Chloro-3-methylphenol | 21 | U | 21 | 150 |
| 4-Chloroaniline | 17 | U | 17 | 150 |
| 4-Chlorophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Nitroaniline | 26 | U | 26 | 200 |
| 4-Nitrophenol | 17 | U | 17 | 330 |
| Acenaphthene | 0.76 | U | 0.76 | 6.7 |
| Acenaphthylene | 0.35 | U | 0.35 | 6.7 |
| Acetophenone | 9.2 | U | 9.2 | 100 |
| Anthracene | 0.78 | U | 0.78 | 6.7 |
| Atrazine | 9.1 | U | 9.1 | 200 |
| Benzaldehyde | 12 | U | 12 | 100 |
| Benzo[a]anthracene | 0.63 | U | 0.63 | 6.7 |
| Benzo[a]pyrene | 0.64 | U | 0.64 | 6.7 |
| Benzo[b]fluoranthene | 0.59 | U | 0.59 | 6.7 |
| Benzo[g,h,i]perylene | 0.35 | U | 0.35 | 6.7 |
| Benzo[k]fluoranthene | 0.68 | U | 0.68 | 6.7 |
| Bis(2-chloroethoxy)methane | 22 | U | 22 | 100 |
| Bis(2-chloroethyl)ether | 2.0 | U | 2.0 | 100 |
| Bis(2-ethylhexyl) phthalate | 61.2 | J | 19 | 70 |
| Butyl benzyl phthalate | 10 | U | 10 | 70 |
| Caprolactam | 37 | U | 37 | 330 |
| Carbazole | 27 | U | 27 | 50 |
| Chrysene | 1.1 | U | 1.1 | 6.7 |
| Dibenz(a,h)anthracene | 0.66 | U | 0.66 | 6.7 |
| Dibenzofuran | 0.66 | U | 0.66 | 50 |
| Diethyl phthalate | 16 | U | 16 | 70 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-206908**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206908/21-A | Analysis Batch: | 240-207287 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206908 | Lab File ID: | 51118003.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/18/2015 0851 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Dimethyl phthalate | 17 | U | 17 | 70 |
| Di-n-butyl phthalate | 15 | U | 15 | 70 |
| Di-n-octyl phthalate | 7.9 | U | 7.9 | 70 |
| Fluoranthene | 0.55 | U | 0.55 | 6.7 |
| Fluorene | 0.53 | U | 0.53 | 6.7 |
| Hexachlorobenzene | 2.1 | U | 2.1 | 6.7 |
| Hexachlorobutadiene | 5.6 | U | 5.6 | 50 |
| Hexachlorocyclopentadiene | 8.1 | U | 8.1 | 330 |
| Hexachloroethane | 9.0 | U | 9.0 | 50 |
| Indeno[1,2,3-cd]pyrene | 0.35 | U | 0.35 | 6.7 |
| Isophorone | 13 | U | 13 | 50 |
| N-Nitrosodi-n-propylamine | 6.3 | U | 6.3 | 50 |
| N-Nitrosodiphenylamine | 21 | U | 21 | 50 |
| Naphthalene | 0.82 | U | 0.82 | 6.7 |
| Nitrobenzene | 2.2 | U | 2.2 | 100 |
| Pentachlorophenol | 9.1 | U | 9.1 | 150 |
| Phenanthrene | 0.73 | U | 0.73 | 6.7 |
| Phenol | 7.3 | U | 7.3 | 50 |
| Pyrene | 0.44 | U | 0.44 | 6.7 |
| 3 & 4 Methylphenol | 20 | U | 20 | 400 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 74 | 36 - 110 |
| Phenol-d5 (Surr) | 73 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 67 | 20 - 110 |
| 2-Fluorophenol (Surr) | 70 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 70 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 30 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-206908**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206908/22-A | Analysis Batch: | 240-207287 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206908 | Lab File ID: | 51118004.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/18/2015 0915 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 442 | 66 | 35 - 110 | |
| bis (2-chloroisopropyl) ether | 667 | 446 | 67 | 29 - 110 | |
| 2,4,5-Trichlorophenol | 667 | 387 | 58 | 25 - 110 | |
| 2,4,6-Trichlorophenol | 667 | 286 | 43 | 12 - 110 | |
| 2,4-Dichlorophenol | 667 | 461 | 69 | 39 - 110 | |
| 2,4-Dimethylphenol | 667 | 467 | 70 | 29 - 110 | |
| 2,4-Dinitrophenol | 1330 | 611 | 46 | 10 - 110 | |
| 2,4-Dinitrotoluene | 667 | 488 | 73 | 48 - 110 | |
| 2,6-Dinitrotoluene | 667 | 480 | 72 | 45 - 110 | |
| 2-Chloronaphthalene | 667 | 463 | 69 | 32 - 110 | |
| 2-Chlorophenol | 667 | 445 | 67 | 37 - 110 | |
| 2-Methylnaphthalene | 667 | 458 | 69 | 36 - 110 | |
| 2-Methylphenol | 667 | 453 | 68 | 41 - 110 | |
| 2-Nitroaniline | 667 | 485 | 73 | 45 - 110 | |
| 2-Nitrophenol | 667 | 461 | 69 | 34 - 110 | |
| 3,3'-Dichlorobenzidine | 1330 | 732 | 55 | 28 - 110 | |
| 3-Nitroaniline | 667 | 432 | 65 | 44 - 110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 779 | 58 | 10 - 110 | |
| 4-Bromophenyl phenyl ether | 667 | 488 | 73 | 39 - 110 | |
| 4-Chloro-3-methylphenol | 667 | 474 | 71 | 48 - 110 | |
| 4-Chloroaniline | 667 | 349 | 52 | 30 - 110 | |
| 4-Chlorophenyl phenyl ether | 667 | 467 | 70 | 40 - 110 | |
| 4-Nitroaniline | 667 | 503 | 75 | 48 - 110 | |
| 4-Nitrophenol | 1330 | 881 | 66 | 28 - 110 | |
| Acenaphthene | 667 | 464 | 70 | 38 - 110 | |
| Acenaphthylene | 667 | 462 | 69 | 40 - 110 | |
| Acetophenone | 667 | 438 | 66 | 40 - 110 | |
| Anthracene | 667 | 471 | 71 | 48 - 110 | |
| Atrazine | 1330 | 971 | 73 | 54 - 120 | |
| Benzaldehyde | 1330 | 909 | 68 | 32 - 110 | |
| Benzo[a]anthracene | 667 | 466 | 70 | 50 - 110 | |
| Benzo[a]pyrene | 667 | 466 | 70 | 44 - 110 | |
| Benzo[b]fluoranthene | 667 | 452 | 68 | 43 - 110 | |
| Benzo[g,h,i]perylene | 667 | 476 | 71 | 51 - 110 | |
| Benzo[k]fluoranthene | 667 | 469 | 70 | 38 - 105 | |
| Bis(2-chloroethoxy)methane | 667 | 486 | 73 | 32 - 110 | |
| Bis(2-chloroethyl)ether | 667 | 453 | 68 | 34 - 110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 704 | 106 | 50 - 110 | |
| Butyl benzyl phthalate | 667 | 489 | 73 | 51 - 110 | |
| Caprolactam | 1330 | 1050 | 79 | 44 - 114 | |
| Carbazole | 667 | 488 | 73 | 50 - 110 | |
| Chrysene | 667 | 495 | 74 | 50 - 110 | |
| Dibenz(a,h)anthracene | 667 | 469 | 70 | 51 - 110 | |
| Dibenzofuran | 667 | 464 | 70 | 43 - 110 | |
| Diethyl phthalate | 667 | 490 | 73 | 52 - 110 | |
| Dimethyl phthalate | 667 | 479 | 72 | 50 - 110 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Lab Control Sample - Batch: 240-206908**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206908/22-A | Analysis Batch: | 240-207287 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206908 | Lab File ID: | 51118004.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/18/2015 0915 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | 667 | 489 | 73 | 51 - 110 | |
| Di-n-octyl phthalate | 667 | 445 | 67 | 48 - 110 | |
| Fluoranthene | 667 | 473 | 71 | 51 - 110 | |
| Fluorene | 667 | 454 | 68 | 46 - 110 | |
| Hexachlorobenzene | 667 | 477 | 72 | 43 - 110 | |
| Hexachlorobutadiene | 667 | 447 | 67 | 29 - 110 | |
| Hexachlorocyclopentadiene | 667 | 397 | 60 | 12 - 110 | |
| Hexachloroethane | 667 | 418 | 63 | 30 - 110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 477 | 72 | 50 - 110 | |
| Isophorone | 667 | 478 | 72 | 36 - 110 | |
| N-Nitrosodi-n-propylamine | 667 | 451 | 68 | 38 - 110 | |
| N-Nitrosodiphenylamine | 1330 | 946 | 71 | 46 - 110 | |
| Naphthalene | 667 | 462 | 69 | 36 - 110 | |
| Nitrobenzene | 667 | 467 | 70 | 32 - 110 | |
| Pentachlorophenol | 1330 | 669 | 50 | 10 - 110 | |
| Phenanthrene | 667 | 471 | 71 | 49 - 110 | |
| Phenol | 667 | 456 | 68 | 38 - 110 | |
| Pyrene | 667 | 475 | 71 | 49 - 110 | |
| 3 & 4 Methylphenol | 667 | 448 | 67 | 40 - 110 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 71 | 36 - 110 |
| Phenol-d5 (Surr) | 71 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 70 | 20 - 110 |
| 2-Fluorophenol (Surr) | 70 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 70 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 42 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206908**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57930-5 | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206908 | Lab File ID: | 51119015.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30.06 g |
| Analysis Date: | 11/19/2015 1500 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57930-5 | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206908 | Lab File ID: | 51119016.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.87 g |
| Analysis Date: | 11/19/2015 1524 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1'-Biphenyl | 70 | 67 | 32 - 110 | 3 | 32 | | |
| bis (2-chloroisopropyl) ether | 71 | 68 | 11 - 110 | 5 | 42 | | |
| 2,4,5-Trichlorophenol | 70 | 70 | 10 - 117 | 0 | 99 | | |
| 2,4,6-Trichlorophenol | 69 | 64 | 10 - 110 | 7 | 38 | | |
| 2,4-Dichlorophenol | 72 | 72 | 10 - 110 | 0 | 34 | | |
| 2,4-Dimethylphenol | 71 | 68 | 10 - 110 | 4 | 31 | | |
| 2,4-Dinitrophenol | 48 | 47 | 10 - 110 | 1 | 99 | | |
| 2,4-Dinitrotoluene | 74 | 75 | 32 - 110 | 2 | 30 | | |
| 2,6-Dinitrotoluene | 72 | 72 | 35 - 110 | 0 | 30 | | |
| 2-Chloronaphthalene | 70 | 69 | 28 - 110 | 0 | 30 | | |
| 2-Chlorophenol | 75 | 69 | 10 - 110 | 7 | 47 | | |
| 2-Methylnaphthalene | 72 | 72 | 10 - 133 | 1 | 42 | | |
| 2-Methylphenol | 75 | 70 | 24 - 110 | 6 | 51 | | |
| 2-Nitroaniline | 71 | 72 | 39 - 110 | 2 | 31 | | |
| 2-Nitrophenol | 70 | 70 | 10 - 110 | 0 | 49 | | |
| 3,3'-Dichlorobenzidine | 54 | 57 | 10 - 110 | 5 | 56 | | |
| 3-Nitroaniline | 63 | 61 | 10 - 110 | 2 | 30 | | |
| 4,6-Dinitro-2-methylphenol | 54 | 55 | 10 - 110 | 2 | 55 | | |
| 4-Bromophenyl phenyl ether | 76 | 75 | 33 - 110 | 1 | 30 | | |
| 4-Chloro-3-methylphenol | 76 | 73 | 25 - 110 | 3 | 54 | | |
| 4-Chloroaniline | 51 | 53 | 10 - 110 | 4 | 36 | | |
| 4-Chlorophenyl phenyl ether | 72 | 72 | 32 - 110 | 1 | 30 | | |
| 4-Nitroaniline | 65 | 66 | 10 - 110 | 3 | 48 | | |
| 4-Nitrophenol | 65 | 64 | 10 - 113 | 1 | 49 | | |
| Acenaphthene | 71 | 69 | 22 - 110 | 3 | 99 | | |
| Acenaphthylene | 71 | 70 | 24 - 110 | 2 | 99 | | |
| Acetophenone | 71 | 68 | 31 - 110 | 4 | 43 | | |
| Anthracene | 70 | 71 | 20 - 110 | 2 | 99 | | |
| Atrazine | 73 | 75 | 45 - 118 | 4 | 30 | | |
| Benzaldehyde | 74 | 69 | 23 - 110 | 5 | 42 | | |
| Benzo[a]anthracene | 72 | 73 | 10 - 122 | 2 | 99 | | |
| Benzo[a]pyrene | 72 | 74 | 10 - 110 | 4 | 99 | | |
| Benzo[b]fluoranthene | 71 | 73 | 12 - 118 | 3 | 99 | | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206908**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57930-5 | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206908 | Lab File ID: | 51119015.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30.06 g |
| Analysis Date: | 11/19/2015 1500 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57930-5 | Analysis Batch: | 240-207541 | Instrument ID: | A4HP10 |
| Client Matrix: | Solid | Prep Batch: | 240-206908 | Lab File ID: | 51119016.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.87 g |
| Analysis Date: | 11/19/2015 1524 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 11/16/2015 0920 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Benzo[g,h,i]perylene | 71 | 72 | 10 - 117 | 2 | 99 | | |
| Benzo[k]fluoranthene | 73 | 76 | 10 - 121 | 4 | 99 | | |
| Bis(2-chloroethoxy)methane | 77 | 75 | 26 - 110 | 2 | 37 | | |
| Bis(2-chloroethyl)ether | 75 | 71 | 21 - 110 | 5 | 55 | | |
| Bis(2-ethylhexyl) phthalate | 75 | 76 | 40 - 110 | 2 | 30 | | |
| Butyl benzyl phthalate | 77 | 78 | 44 - 110 | 3 | 30 | | |
| Caprolactam | 83 | 83 | 10 - 134 | 1 | 32 | | |
| Carbazole | 68 | 69 | 34 - 110 | 1 | 30 | | |
| Chrysene | 75 | 77 | 10 - 125 | 4 | 99 | | |
| Dibenz(a,h)anthracene | 72 | 73 | 14 - 113 | 2 | 99 | | |
| Dibenzofuran | 71 | 70 | 29 - 110 | 2 | 30 | | |
| Diethyl phthalate | 75 | 75 | 42 - 110 | 0 | 30 | | |
| Dimethyl phthalate | 73 | 73 | 41 - 110 | 0 | 30 | | |
| Di-n-butyl phthalate | 74 | 75 | 43 - 110 | 1 | 30 | | |
| Di-n-octyl phthalate | 73 | 75 | 24 - 119 | 4 | 30 | | |
| Fluoranthene | 73 | 73 | 10 - 110 | 1 | 99 | | |
| Fluorene | 70 | 69 | 23 - 110 | 1 | 99 | | |
| Hexachlorobenzene | 70 | 71 | 34 - 110 | 2 | 30 | | |
| Hexachlorobutadiene | 69 | 69 | 25 - 110 | 1 | 34 | | |
| Hexachlorocyclopentadiene | 40 | 30 | 10 - 110 | 28 | 79 | J | J |
| Hexachloroethane | 66 | 61 | 12 - 110 | 7 | 50 | | |
| Indeno[1,2,3-cd]pyrene | 71 | 73 | 10 - 114 | 4 | 99 | | |
| Isophorone | 76 | 74 | 29 - 110 | 2 | 38 | | |
| N-Nitrosodi-n-propylamine | 72 | 68 | 26 - 110 | 5 | 42 | | |
| N-Nitrosodiphenylamine | 72 | 71 | 22 - 110 | 1 | 30 | | |
| Naphthalene | 72 | 71 | 10 - 111 | 1 | 99 | | |
| Nitrobenzene | 73 | 71 | 23 - 110 | 2 | 41 | | |
| Pentachlorophenol | 55 | 50 | 10 - 110 | 8 | 50 | | |
| Phenanthrene | 71 | 72 | 10 - 166 | 1 | 99 | | |
| Phenol | 74 | 70 | 17 - 110 | 5 | 53 | | |
| Pyrene | 74 | 74 | 10 - 147 | 1 | 99 | | |
| 3 & 4 Methylphenol | 74 | 69 | 25 - 110 | 6 | 50 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 76 | 75 | 36 - 110 |
| Phenol-d5 (Surr) | 78 | 74 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 73 | 71 | 20 - 110 |
| 2-Fluorophenol (Surr) | 78 | 73 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 74 | 72 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 72 | 64 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206908**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57930-5 | Units: ug/Kg | | MSD Lab Sample ID: | 240-57930-5 |
| Client Matrix: | Solid | | | Client Matrix: | Solid |
| Dilution: | 1.0 | | | Dilution: | 1.0 |
| Analysis Date: | 11/19/2015 1500 | | | Analysis Date: | 11/19/2015 1524 |
| Prep Date: | 11/16/2015 0920 | | | Prep Date: | 11/16/2015 0920 |
| Leach Date: | N/A | | | Leach Date: | N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 4.1 | U | 786 | 791 | 547 | 534 |
| bis (2-chloroisopropyl) ether | 11 | U | 786 | 791 | 561 | 534 |
| 2,4,5-Trichlorophenol | 29 | U | 786 | 791 | 554 | 553 |
| 2,4,6-Trichlorophenol | 10 | U | 786 | 791 | 543 | 506 |
| 2,4-Dichlorophenol | 23 | U | 786 | 791 | 567 | 568 |
| 2,4-Dimethylphenol | 23 | U | 786 | 791 | 554 | 535 |
| 2,4-Dinitrophenol | 25 | U | 1570 | 1580 | 752 | 741 |
| 2,4-Dinitrotoluene | 20 | U | 786 | 791 | 582 | 593 |
| 2,6-Dinitrotoluene | 25 | U | 786 | 791 | 567 | 567 |
| 2-Chloronaphthalene | 0.53 | U | 786 | 791 | 547 | 545 |
| 2-Chlorophenol | 9.6 | U | 786 | 791 | 587 | 546 |
| 2-Methylnaphthalene | 0.59 | U | 786 | 791 | 564 | 568 |
| 2-Methylphenol | 13 | U | 786 | 791 | 587 | 554 |
| 2-Nitroaniline | 11 | U | 786 | 791 | 558 | 568 |
| 2-Nitrophenol | 9.7 | U | 786 | 791 | 553 | 555 |
| 3,3'-Dichlorobenzidine | 21 | U | 1570 | 1580 | 856 | 899 |
| 3-Nitroaniline | 19 | U | 786 | 791 | 491 | 483 |
| 4,6-Dinitro-2-methylphenol | 11 | U | 1570 | 1580 | 856 | 875 |
| 4-Bromophenyl phenyl ether | 15 | U | 786 | 791 | 597 | 593 |
| 4-Chloro-3-methylphenol | 25 | U | 786 | 791 | 598 | 579 |
| 4-Chloroaniline | 20 | U | 786 | 791 | 401 | 419 |
| 4-Chlorophenyl phenyl ether | 15 | U | 786 | 791 | 567 | 571 |
| 4-Nitroaniline | 30 | U | 786 | 791 | 510 | 524 |
| 4-Nitrophenol | 20 | U | 1570 | 1580 | 1030 | 1020 |
| Acenaphthene | 0.89 | U | 786 | 791 | 557 | 542 |
| Acenaphthylene | 0.41 | U | 786 | 791 | 562 | 552 |
| Acetophenone | 11 | U | 786 | 791 | 559 | 538 |
| Anthracene | 0.91 | U | 786 | 791 | 552 | 564 |
| Atrazine | 11 | U | 1570 | 1580 | 1150 | 1190 |
| Benzaldehyde | 14 | U | 1570 | 1580 | 1160 | 1100 |
| Benzo[a]anthracene | 0.74 | U | 786 | 791 | 567 | 580 |
| Benzo[a]pyrene | 0.75 | U | 786 | 791 | 564 | 588 |
| Benzo[b]fluoranthene | 0.69 | U | 786 | 791 | 560 | 575 |
| Benzo[g,h,i]perylene | 0.41 | U | 786 | 791 | 561 | 571 |
| Benzo[k]fluoranthene | 0.80 | U | 786 | 791 | 576 | 598 |
| Bis(2-chloroethoxy)methane | 26 | U | 786 | 791 | 605 | 594 |
| Bis(2-chloroethyl)ether | 2.3 | U | 786 | 791 | 593 | 564 |
| Bis(2-ethylhexyl) phthalate | 25 | J | 786 | 791 | 613 | 627 |
| Butyl benzyl phthalate | 12 | U | 786 | 791 | 601 | 617 |
| Caprolactam | 43 | U | 1570 | 1580 | 1300 | 1320 |
| Carbazole | 32 | U | 786 | 791 | 537 | 545 |
| Chrysene | 1.3 | U | 786 | 791 | 586 | 608 |
| Dibenz(a,h)anthracene | 0.77 | U | 786 | 791 | 564 | 575 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206908**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57930-5 | Units: ug/Kg | MSD Lab Sample ID: | 240-57930-5 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 11/19/2015 1500 | | Analysis Date: | 11/19/2015 1524 | |
| Prep Date: | 11/16/2015 0920 | | Prep Date: | 11/16/2015 0920 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| Dibenzofuran | 0.77 | U | 786 | 791 | 562 | | 552 | |
| Diethyl phthalate | 19 | U | 786 | 791 | 591 | | 590 | |
| Dimethyl phthalate | 20 | U | 786 | 791 | 575 | | 576 | |
| Di-n-butyl phthalate | 18 | U | 786 | 791 | 585 | | 593 | |
| Di-n-octyl phthalate | 9.3 | U | 786 | 791 | 570 | | 592 | |
| Fluoranthene | 0.64 | U | 786 | 791 | 574 | | 578 | |
| Fluorene | 0.62 | U | 786 | 791 | 553 | | 545 | |
| Hexachlorobenzene | 2.5 | U | 786 | 791 | 551 | | 562 | |
| Hexachlorobutadiene | 6.6 | U | 786 | 791 | 543 | | 548 | |
| Hexachlorocyclopentadiene | 9.5 | U | 786 | 791 | 315 | J | 238 | J |
| Hexachloroethane | 11 | U | 786 | 791 | 519 | | 485 | |
| Indeno[1,2,3-cd]pyrene | 0.41 | U | 786 | 791 | 558 | | 579 | |
| Isophorone | 15 | U | 786 | 791 | 596 | | 586 | |
| N-Nitrosodi-n-propylamine | 7.4 | U | 786 | 791 | 570 | | 540 | |
| N-Nitrosodiphenylamine | 25 | U | 1570 | 1580 | 1140 | | 1130 | |
| Naphthalene | 0.96 | U | 786 | 791 | 563 | | 559 | |
| Nitrobenzene | 2.6 | U | 786 | 791 | 571 | | 562 | |
| Pentachlorophenol | 11 | U | 1570 | 1580 | 863 | | 793 | |
| Phenanthrene | 0.86 | U | 786 | 791 | 560 | | 567 | |
| Phenol | 8.6 | U | 786 | 791 | 583 | | 554 | |
| Pyrene | 0.52 | U | 786 | 791 | 579 | | 586 | |
| 3 & 4 Methylphenol | 23 | U | 786 | 791 | 578 | | 544 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-206878**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206878/1-A | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Client Matrix: | Solid | Prep Batch: | 240-206878 | Lab File ID: | Y0111609.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/16/2015 1010 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 46 | U | 46 | 100 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 71 | 40 - 139 |

**Lab Control Sample - Batch: 240-206878**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206878/2-A | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Client Matrix: | Solid | Prep Batch: | 240-206878 | Lab File ID: | Y0111610.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/16/2015 1045 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 651 | 81 | 74 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 72 | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206878**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57930-11 | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Client Matrix: | Solid | Prep Batch: | 240-206878 | Lab File ID: | Y0111623.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 0.47 g |
| Analysis Date: | 11/16/2015 1814 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57930-11 | Analysis Batch: | 240-206833 | Instrument ID: | YPID |
| Client Matrix: | Solid | Prep Batch: | 240-206878 | Lab File ID: | Y0111624.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 0.51 g |
| Analysis Date: | 11/16/2015 1849 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/16/2015 0812 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Gasoline Range Organics [C6 - C10] | 80 | 78 | 10 - 136 | 9 | 40 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 87 | 73 | 40 - 139 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206878**

**Method: 8015C**
**Preparation: 5030A**

| | | | |
|---|---|---|---|
| MS Lab Sample ID: | 240-57930-11 | Units: ug/Kg | MSD Lab Sample ID: 240-57930-11 |
| Client Matrix: | Solid | | Client Matrix: Solid |
| Dilution: | 1.0 | | Dilution: 1.0 |
| Analysis Date: | 11/16/2015 1814 | | Analysis Date: 11/16/2015 1849 |
| Prep Date: | 11/16/2015 0812 | | Prep Date: 11/16/2015 0812 |
| Leach Date: | N/A | | Leach Date: N/A |

| Analyte | Sample Result/Qual | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 1800 | 10100 | 9290 | 9930 | 9080 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-207078**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207078/1-A | Analysis Batch: | 240-207037 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-207078 | Lab File ID: | A0111705.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/17/2015 0924 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 0748 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 47.9 | J | 46 | 100 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 79 | 40 - 139 |

**Lab Control Sample - Batch: 240-207078**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207078/2-A | Analysis Batch: | 240-207037 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-207078 | Lab File ID: | A0111706.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/17/2015 1001 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 0748 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 889 | 111 | 74 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 101 | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207078**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57774-C-8-J MS | Analysis Batch: | 240-207037 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-207078 | Lab File ID: | A0111715.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 4.92 g |
| Analysis Date: | 11/17/2015 1535 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 0748 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57774-C-8-K MSD | Analysis Batch: | 240-207037 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-207078 | Lab File ID: | A0111716.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 4.94 g |
| Analysis Date: | 11/17/2015 1612 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 0748 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Gasoline Range Organics [C6 - C10] | 101 | 101 | 10 - 136 | 0 | 40 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 100 | 100 | 40 - 139 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207078**

**Method: 8015C**
**Preparation: 5030A**

| | | | |
|---|---|---|---|
| MS Lab Sample ID: | 240-57774-C-8-J MS | Units: | ug/Kg |
| Client Matrix: | Solid | | |
| Dilution: | 1.0 | | |
| Analysis Date: | 11/17/2015 1535 | | |
| Prep Date: | 11/17/2015 0748 | | |
| Leach Date: | N/A | | |

| | |
|---|---|
| MSD Lab Sample ID: | 240-57774-C-8-K MSD |
| Client Matrix: | Solid |
| Dilution: | 1.0 |
| Analysis Date: | 11/17/2015 1612 |
| Prep Date: | 11/17/2015 0748 |
| Leach Date: | N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 52 | U | 899 | 896 | 906 | 902 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-207274**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207274/1-A | Analysis Batch: | 240-207246 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-207274 | Lab File ID: | A0111817.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/18/2015 1548 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 0657 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 46 | U | 46 | 100 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 78 | 40 - 139 |

**Lab Control Sample - Batch: 240-207274**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207274/2-A | Analysis Batch: | 240-207246 | Instrument ID: | AFID |
| Client Matrix: | Solid | Prep Batch: | 240-207274 | Lab File ID: | A0111818.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 11/18/2015 1626 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/18/2015 0657 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 836 | 104 | 74 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 101 | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-206884**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206884/23-A | Analysis Batch: | 240-207234 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-206884 | Lab File ID: | P5B11708.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/17/2015 2208 | Units: | mg/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/16/2015 0833 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | 6.4 | U | 6.4 | 17 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| o-Terphenyl (Surr) | 77 | 40 - 160 |

**Lab Control Sample - Batch: 240-206884**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206884/24-A | Analysis Batch: | 240-207234 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-206884 | Lab File ID: | P5B11709.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/17/2015 2240 | Units: | mg/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/16/2015 0833 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Diesel | 84.0 | 66.9 | 80 | 52 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| o-Terphenyl (Surr) | 82 | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-207121**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-207121/7-A | Analysis Batch: | 240-207465 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-207121 | Lab File ID: | P5B11815.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/19/2015 0129 | Units: | mg/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1002 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | 6.4 | U | 6.4 | 17 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| o-Terphenyl (Surr) | 75 | 40 - 160 |

**Lab Control Sample - Batch: 240-207121**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-207121/8-A | Analysis Batch: | 240-207465 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-207121 | Lab File ID: | P5B11816.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/19/2015 0201 | Units: | mg/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1002 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Diesel | 84.0 | 83.2 | 99 | 52 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| o-Terphenyl (Surr) | 103 | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207121**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57930-11 | Analysis Batch: | 240-207465 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-207121 | Lab File ID: | P5B11821.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.91 g |
| Analysis Date: | 11/19/2015 0438 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1002 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57930-11 | Analysis Batch: | 240-207465 | Instrument ID: | A2HP5R |
| Client Matrix: | Solid | Prep Batch: | 240-207121 | Lab File ID: | P5B11822.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30.50 g |
| Analysis Date: | 11/19/2015 0510 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 11/17/2015 1002 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Diesel | 74 | 79 | 10 - 124 | 3 | 40 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 93 | 93 | 40 - 160 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-207121**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57930-11 | Units: mg/Kg | | MSD Lab Sample ID: | 240-57930-11 |
| Client Matrix: | Solid | | | Client Matrix: | Solid |
| Dilution: | 1.0 | | | Dilution: | 1.0 |
| Analysis Date: | 11/19/2015 0438 | | | Analysis Date: | 11/19/2015 0510 |
| Prep Date: | 11/17/2015 1002 | | | Prep Date: | 11/17/2015 1002 |
| Leach Date: | N/A | | | Leach Date: | N/A |

| Analyte | Sample Result/Qual | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|
| Diesel | 34 | 99.9 | 97.9 | 108 | 112 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Method Blank - Batch: 240-206917**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-206917/21-A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Client Matrix: | Solid | Prep Batch: | 240-206917 | Lab File ID: | P1200062.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/19/2015 0103 | Units: | ug/Kg | Final Weight/Volume: | 10 mL |
| Prep Date: | 11/16/2015 0932 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Aroclor-1016 | 12 | U | 12 | 33 |
| Aroclor-1221 | 16 | U | 16 | 33 |
| Aroclor-1232 | 20 | U | 20 | 33 |
| Aroclor-1242 | 11 | U | 11 | 33 |
| Aroclor-1248 | 8.0 | U | 8.0 | 33 |
| Aroclor-1254 | 14 | U | 14 | 33 |
| Aroclor-1260 | 9.0 | U | 9.0 | 33 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Tetrachloro-m-xylene | 74 | 13 - 134 |
| DCB Decachlorobiphenyl | 79 | 10 - 155 |

**Lab Control Sample - Batch: 240-206917**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-206917/22-A | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Client Matrix: | Solid | Prep Batch: | 240-206917 | Lab File ID: | P1200063.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 11/19/2015 0121 | Units: | ug/Kg | Final Weight/Volume: | 10 mL |
| Prep Date: | 11/16/2015 0932 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Aroclor-1016 | 333 | 251 | 75 | 51 - 120 | |
| Aroclor-1260 | 334 | 253 | 76 | 48 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Tetrachloro-m-xylene | 75 | 13 - 134 |
| DCB Decachlorobiphenyl | 77 | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206917**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-57880-E-1-B MS | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Client Matrix: | Solid | Prep Batch: | 240-206917 | Lab File ID: | P1200066.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.95 g |
| Analysis Date: | 11/19/2015 0216 | | | Final Weight/Volume: | 10 mL |
| Prep Date: | 11/16/2015 0932 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-57880-E-1-C MSD | Analysis Batch: | 240-207364 | Instrument ID: | A2HP12 |
| Client Matrix: | Solid | Prep Batch: | 240-206917 | Lab File ID: | P1200067.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.75 g |
| Analysis Date: | 11/19/2015 0234 | | | Final Weight/Volume: | 10 mL |
| Prep Date: | 11/16/2015 0932 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Aroclor-1016 | 117 | 143 | 17 - 141 | 18 | 56 | | F1 |
| Aroclor-1260 | 44 | 55 | 19 - 137 | 21 | 55 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 75 | 90 | 13 - 134 |
| DCB Decachlorobiphenyl | 94 | 87 | 10 - 155 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-206917**

**Method: 8082A**
**Preparation: 3540C**

| | | | |
|---|---|---|---|
| MS Lab Sample ID: | 240-57880-E-1-B MS | Units: ug/Kg | MSD Lab Sample ID: 240-57880-E-1-C MSD |
| Client Matrix: | Solid | | Client Matrix: Solid |
| Dilution: | 1.0 | | Dilution: 1.0 |
| Analysis Date: | 11/19/2015 0216 | | Analysis Date: 11/19/2015 0234 |
| Prep Date: | 11/16/2015 0932 | | Prep Date: 11/16/2015 0932 |
| Leach Date: | N/A | | Leach Date: N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|
| Aroclor-1016 | 16 | U | 443 | 446 | 518 | 638 | F1 |
| Aroclor-1260 | 71 | | 444 | 447 | 266 | 331 | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

**Duplicate - Batch: 240-207016**

**Method: Moisture**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | 240-57930-9 | Analysis Batch: | 240-207016 | Instrument ID: | No Equipment Assigned |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | N/A |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | |
| Analysis Date: | 11/16/2015 1541 | Units: | % | Final Weight/Volume: | |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Sample Result/Qual | Result | RPD | Limit | Qual |
|---|---|---|---|---|---|
| Percent Solids | 83 | 83 | 0.2 | 20 | |
| Percent Moisture | 17 | 17 | 0.9 | 20 | |

# DATA REPORTING QUALIFIERS

Client: Tetra Tech, Inc. Job Number: 240-57930-1

| Lab Section | Qualifier | Description |
|---|---|---|
| GC/MS VOA | | |
| | B | Compound was found in the blank and sample. |
| | ^ | ICV,CCV,ICB,CCB, ISA, ISB, CRI, CRA, DLCK or MRL standard: Instrument related QC is outside acceptance limits. |
| | U | Indicates the analyte was analyzed for but not detected. |
| | * | ISTD response or retention time outside acceptable limits |
| | * | LCS or LCSD is outside acceptance limits. |
| | F1 | MS and/or MSD Recovery is outside acceptance limits. |
| | E | Result exceeded calibration range. |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| GC/MS Semi VOA | | |
| | B | Compound was found in the blank and sample. |
| | U | Indicates the analyte was analyzed for but not detected. |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| GC VOA | | |
| | B | Compound was found in the blank and sample. |
| | U | Indicates the analyte was analyzed for but not detected. |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| GC Semi VOA | | |
| | U | Indicates the analyte was analyzed for but not detected. |
| | F1 | MS and/or MSD Recovery is outside acceptance limits. |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC/MS VOA** | | | | | |
| **Prep Batch: 240-206968** | | | | | |
| 240-57921-A-4-J MS | Matrix Spike | T | Solid | | |
| 240-57921-A-4-K MSD | Matrix Spike Duplicate | T | Solid | | |
| **Analysis Batch:240-207096** | | | | | |
| LCS 240-207096/4 | Lab Control Sample | T | Water | 8260C | |
| MB 240-207096/6 | Method Blank | T | Water | 8260C | |
| 200-30727-B-1 MS | Matrix Spike | T | Water | 8260C | |
| 200-30727-B-1 MSD | Matrix Spike Duplicate | T | Water | 8260C | |
| 240-57930-12 | TB-111315 | T | Water | 8260C | |
| **Analysis Batch:240-207343** | | | | | |
| LCS 240-207343/5 | Lab Control Sample | T | Solid | 8260C | |
| LCS 240-207343/7 | Lab Control Sample | T | Solid | 8260C | |
| MB 240-207343/6 | Method Blank | T | Solid | 8260C | |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 8260C | 240-207393 |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 8260C | 240-207393 |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 8260C | 240-207393 |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 8260C | 240-207393 |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 8260C | 240-207393 |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 8260C | 240-207393 |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 8260C | 240-207393 |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 8260C | 240-207393 |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 8260C | 240-207393 |
| **Prep Batch: 240-207393** | | | | | |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 5030C | |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 5030C | |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 5030C | |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 5030C | |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 5030C | |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 5030C | |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 5030C | |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 5030C | |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 5030C | |
| **Analysis Batch:240-207560** | | | | | |
| LCS 240-207560/5 | Lab Control Sample | T | Solid | 8260C | |
| LCS 240-207560/7 | Lab Control Sample | T | Solid | 8260C | |
| MB 240-207560/6 | Method Blank | T | Solid | 8260C | |
| 240-57921-A-4-J MS | Matrix Spike | T | Solid | 8260C | 240-207613 |
| 240-57921-A-4-K MSD | Matrix Spike Duplicate | T | Solid | 8260C | 240-207613 |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 8260C | 240-207613 |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 8260C | 240-207613 |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC/MS VOA** | | | | | |
| **Prep Batch: 240-207613** | | | | | |
| 240-57921-A-4-J MS | Matrix Spike | T | Solid | 5030C | 240-206968 |
| 240-57921-A-4-K MSD | Matrix Spike Duplicate | T | Solid | 5030C | 240-206968 |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 5030C | |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 5030C | |
| **Prep Batch: 240-207709** | | | | | |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 5030C | |
| 240-57972-C-3-C MS | Matrix Spike | T | Solid | 5030C | |
| 240-57972-C-3-D MSD | Matrix Spike Duplicate | T | Solid | 5030C | |
| **Analysis Batch:240-207720** | | | | | |
| LCS 240-207720/6 | Lab Control Sample | T | Solid | 8260C | |
| MB 240-207720/8 | Method Blank | T | Solid | 8260C | |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 8260C | 240-207709 |
| 240-57972-C-3-C MS | Matrix Spike | T | Solid | 8260C | 240-207709 |
| 240-57972-C-3-D MSD | Matrix Spike Duplicate | T | Solid | 8260C | 240-207709 |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC/MS Semi VOA** | | | | | |
| **Prep Batch: 240-206887** | | | | | |
| LCS 240-206887/24-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-206887/23-A | Method Blank | T | Solid | 3540C | |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 3540C | |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 3540C | |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 3540C | |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 3540C | |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 3540C | |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 3540C | |
| **Prep Batch: 240-206908** | | | | | |
| LCS 240-206908/22-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-206908/21-A | Method Blank | T | Solid | 3540C | |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 3540C | |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 3540C | |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 3540C | |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 3540C | |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 3540C | |
| 240-57930-5MS | Matrix Spike | T | Solid | 3540C | |
| 240-57930-5MSD | Matrix Spike Duplicate | T | Solid | 3540C | |
| **Analysis Batch:240-207287** | | | | | |
| LCS 240-206908/22-A | Lab Control Sample | T | Solid | 8270D | 240-206908 |
| MB 240-206908/21-A | Method Blank | T | Solid | 8270D | 240-206908 |
| **Analysis Batch:240-207541** | | | | | |
| LCS 240-206887/24-A | Lab Control Sample | T | Solid | 8270D | 240-206887 |
| MB 240-206887/23-A | Method Blank | T | Solid | 8270D | 240-206887 |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 8270D | 240-206908 |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 8270D | 240-206908 |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 8270D | 240-206908 |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 8270D | 240-206908 |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 8270D | 240-206908 |
| 240-57930-5MS | Matrix Spike | T | Solid | 8270D | 240-206908 |
| 240-57930-5MSD | Matrix Spike Duplicate | T | Solid | 8270D | 240-206908 |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 8270D | 240-206887 |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 8270D | 240-206887 |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 8270D | 240-206887 |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 8270D | 240-206887 |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 8270D | 240-206887 |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 8270D | 240-206887 |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC VOA** | | | | | |
| **Analysis Batch:240-206833** | | | | | |
| LCS 240-206878/2-A | Lab Control Sample | T | Solid | 8015C | 240-206878 |
| MB 240-206878/1-A | Method Blank | T | Solid | 8015C | 240-206878 |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 8015C | 240-206878 |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 8015C | 240-206878 |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 8015C | 240-206878 |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 8015C | 240-206878 |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 8015C | 240-206878 |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 8015C | 240-206878 |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 8015C | 240-206878 |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 8015C | 240-206878 |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 8015C | 240-206878 |
| 240-57930-11MS | Matrix Spike | T | Solid | 8015C | 240-206878 |
| 240-57930-11MSD | Matrix Spike Duplicate | T | Solid | 8015C | 240-206878 |
| **Prep Batch: 240-206878** | | | | | |
| LCS 240-206878/2-A | Lab Control Sample | T | Solid | 5030A | |
| MB 240-206878/1-A | Method Blank | T | Solid | 5030A | |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 5030A | |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 5030A | |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 5030A | |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 5030A | |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 5030A | |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 5030A | |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 5030A | |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 5030A | |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 5030A | |
| 240-57930-11MS | Matrix Spike | T | Solid | 5030A | |
| 240-57930-11MSD | Matrix Spike Duplicate | T | Solid | 5030A | |
| **Analysis Batch:240-207037** | | | | | |
| LCS 240-207078/2-A | Lab Control Sample | T | Solid | 8015C | 240-207078 |
| MB 240-207078/1-A | Method Blank | T | Solid | 8015C | 240-207078 |
| 240-57774-C-8-J MS | Matrix Spike | T | Solid | 8015C | 240-207078 |
| 240-57774-C-8-K MSD | Matrix Spike Duplicate | T | Solid | 8015C | 240-207078 |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 8015C | 240-207078 |
| **Prep Batch: 240-207078** | | | | | |
| LCS 240-207078/2-A | Lab Control Sample | T | Solid | 5030A | |
| MB 240-207078/1-A | Method Blank | T | Solid | 5030A | |
| 240-57774-C-8-J MS | Matrix Spike | T | Solid | 5030A | |
| 240-57774-C-8-K MSD | Matrix Spike Duplicate | T | Solid | 5030A | |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 5030A | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC VOA** | | | | | |
| **Analysis Batch:240-207246** | | | | | |
| LCS 240-207274/2-A | Lab Control Sample | T | Solid | 8015C | 240-207274 |
| MB 240-207274/1-A | Method Blank | T | Solid | 8015C | 240-207274 |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 8015C | 240-207274 |
| **Prep Batch: 240-207274** | | | | | |
| LCS 240-207274/2-A | Lab Control Sample | T | Solid | 5030A | |
| MB 240-207274/1-A | Method Blank | T | Solid | 5030A | |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 5030A | |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC Semi VOA** | | | | | |
| **Prep Batch: 240-206884** | | | | | |
| LCS 240-206884/24-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-206884/23-A | Method Blank | T | Solid | 3540C | |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 3540C | |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 3540C | |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 3540C | |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 3540C | |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 3540C | |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 3540C | |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 3540C | |
| **Prep Batch: 240-206917** | | | | | |
| LCS 240-206917/22-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-206917/21-A | Method Blank | T | Solid | 3540C | |
| 240-57880-E-1-B MS | Matrix Spike | T | Solid | 3540C | |
| 240-57880-E-1-C MSD | Matrix Spike Duplicate | T | Solid | 3540C | |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 3540C | |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 3540C | |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 3540C | |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 3540C | |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 3540C | |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 3540C | |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 3540C | |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 3540C | |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 3540C | |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 3540C | |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 3540C | |
| **Prep Batch: 240-207121** | | | | | |
| LCS 240-207121/8-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-207121/7-A | Method Blank | T | Solid | 3540C | |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 3540C | |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 3540C | |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 3540C | |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 3540C | |
| 240-57930-11MS | Matrix Spike | T | Solid | 3540C | |
| 240-57930-11MSD | Matrix Spike Duplicate | T | Solid | 3540C | |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC Semi VOA** | | | | | |
| **Analysis Batch:240-207234** | | | | | |
| LCS 240-206884/24-A | Lab Control Sample | T | Solid | 8015C | 240-206884 |
| MB 240-206884/23-A | Method Blank | T | Solid | 8015C | 240-206884 |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 8015C | 240-206884 |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 8015C | 240-206884 |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 8015C | 240-206884 |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 8015C | 240-206884 |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 8015C | 240-206884 |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 8015C | 240-206884 |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 8015C | 240-206884 |
| **Analysis Batch:240-207364** | | | | | |
| LCS 240-206917/22-A | Lab Control Sample | T | Solid | 8082A | 240-206917 |
| MB 240-206917/21-A | Method Blank | T | Solid | 8082A | 240-206917 |
| 240-57880-E-1-B MS | Matrix Spike | T | Solid | 8082A | 240-206917 |
| 240-57880-E-1-C MSD | Matrix Spike Duplicate | T | Solid | 8082A | 240-206917 |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | 8082A | 240-206917 |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | 8082A | 240-206917 |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | 8082A | 240-206917 |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | 8082A | 240-206917 |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | 8082A | 240-206917 |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | 8082A | 240-206917 |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | 8082A | 240-206917 |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 8082A | 240-206917 |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 8082A | 240-206917 |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 8082A | 240-206917 |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 8082A | 240-206917 |
| **Analysis Batch:240-207465** | | | | | |
| LCS 240-207121/8-A | Lab Control Sample | T | Solid | 8015C | 240-207121 |
| MB 240-207121/7-A | Method Blank | T | Solid | 8015C | 240-207121 |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | 8015C | 240-207121 |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | 8015C | 240-207121 |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | 8015C | 240-207121 |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | 8015C | 240-207121 |
| 240-57930-11MS | Matrix Spike | T | Solid | 8015C | 240-207121 |
| 240-57930-11MSD | Matrix Spike Duplicate | T | Solid | 8015C | 240-207121 |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **General Chemistry** | | | | | |
| **Analysis Batch:240-207016** | | | | | |
| 240-57930-1 | F-VS-UST-P-1 | T | Solid | Moisture | |
| 240-57930-2 | F-VS-UST-P-2 | T | Solid | Moisture | |
| 240-57930-3 | F-VS-UST-P-3 | T | Solid | Moisture | |
| 240-57930-4 | F-VS-UST-P-4 | T | Solid | Moisture | |
| 240-57930-5 | F-VB-UST-P-1 | T | Solid | Moisture | |
| 240-57930-6 | F-VB-UST-P-2 | T | Solid | Moisture | |
| 240-57930-7 | F-VB-UST-P-3 | T | Solid | Moisture | |
| 240-57930-8 | F-VB-UST-P-4 | T | Solid | Moisture | |
| 240-57930-9 | F-VB-UST-P-5 | T | Solid | Moisture | |
| 240-57930-9DU | Duplicate | T | Solid | Moisture | |
| 240-57930-10 | F-VB-UST-P-6 | T | Solid | Moisture | |
| 240-57930-11 | F-VS-UST-P-DUP | T | Solid | Moisture | |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Laboratory Chronicle

**Lab ID: 240-57930-1** **Client ID: F-VS-UST-P-1**

Sample Date/Time: 11/13/2015 10:30 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-1-E | | 240-207343 | 240-207393 | 11/18/2015 12:27 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-1-E | | 240-207343 | 240-207393 | 11/18/2015 18:03 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-1-B | | 240-207541 | 240-206908 | 11/16/2015 09:20 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-1-B | | 240-207541 | 240-206908 | 11/19/2015 19:46 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-1-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-1-A | | 240-206833 | 240-206878 | 11/16/2015 11:18 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-1-A | | 240-207234 | 240-206884 | 11/16/2015 08:33 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-1-A | | 240-207234 | 240-206884 | 11/18/2015 07:32 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-1-C | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-1-C | | 240-207364 | 240-206917 | 11/18/2015 23:51 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-1 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

**Lab ID: 240-57930-2** **Client ID: F-VS-UST-P-2**

Sample Date/Time: 11/13/2015 10:35 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-2-D | | 240-207560 | 240-207613 | 11/19/2015 13:00 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-2-D | | 240-207560 | 240-207613 | 11/19/2015 17:52 | 1 | TAL CAN | SAM |
| P:5030C | 240-57930-B-2-E | | 240-207560 | 240-207613 | 11/19/2015 13:00 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-2-E | | 240-207560 | 240-207613 | 11/19/2015 18:13 | 1 | TAL CAN | SAM |
| P:5030C | 240-57930-B-2-F | | 240-207720 | 240-207709 | 11/20/2015 09:47 | 1 | TAL CAN | TJL2 |
| A:8260C | 240-57930-B-2-F | | 240-207720 | 240-207709 | 11/20/2015 10:06 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-2-B | | 240-207541 | 240-206908 | 11/16/2015 09:20 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-2-B | | 240-207541 | 240-206908 | 11/19/2015 19:23 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-2-B | | 240-207037 | 240-207078 | 11/17/2015 07:48 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-2-B | | 240-207037 | 240-207078 | 11/17/2015 10:39 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-2-A | | 240-207234 | 240-206884 | 11/16/2015 08:33 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-2-A | | 240-207234 | 240-206884 | 11/18/2015 08:04 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-2-C | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-2-C | | 240-207364 | 240-206917 | 11/19/2015 00:09 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-2 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Laboratory Chronicle

**Lab ID: 240-57930-3** **Client ID: F-VS-UST-P-3**

Sample Date/Time: 11/13/2015 10:40 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-3-B | | 240-207343 | 240-207393 | 11/18/2015 12:27 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-3-B | | 240-207343 | 240-207393 | 11/18/2015 18:25 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-3-B | | 240-207541 | 240-206908 | 11/16/2015 09:20 | 2.5 | TAL CAN | YDN |
| A:8270D | 240-57930-D-3-B | | 240-207541 | 240-206908 | 11/19/2015 20:57 | 2.5 | TAL CAN | JMG |
| P:5030A | 240-57930-C-3-B | | 240-207246 | 240-207274 | 11/18/2015 06:57 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-3-B | | 240-207246 | 240-207274 | 11/18/2015 17:04 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-3-A | | 240-207234 | 240-206884 | 11/16/2015 08:33 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-3-A | | 240-207234 | 240-206884 | 11/18/2015 08:35 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-3-C | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-3-C | | 240-207364 | 240-206917 | 11/19/2015 00:27 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-3 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

**Lab ID: 240-57930-4** **Client ID: F-VS-UST-P-4**

Sample Date/Time: 11/13/2015 10:45 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-4-E | | 240-207560 | 240-207613 | 11/19/2015 13:00 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-4-E | | 240-207560 | 240-207613 | 11/19/2015 17:08 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-4-B | | 240-207541 | 240-206908 | 11/16/2015 09:20 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-4-B | | 240-207541 | 240-206908 | 11/19/2015 20:34 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-4-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-4-A | | 240-206833 | 240-206878 | 11/16/2015 13:01 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-4-A | | 240-207234 | 240-206884 | 11/16/2015 08:33 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-4-A | | 240-207234 | 240-206884 | 11/18/2015 10:10 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-4-C | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-4-C | | 240-207364 | 240-206917 | 11/19/2015 00:45 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-4 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Laboratory Chronicle

**Lab ID: 240-57930-5** **Client ID: F-VB-UST-P-1**

Sample Date/Time: 11/13/2015 10:00 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-5-B | | 240-207343 | 240-207393 | 11/18/2015 12:27 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-5-B | | 240-207343 | 240-207393 | 11/18/2015 19:51 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-5-B | | 240-207541 | 240-206908 | 11/16/2015 09:20 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-5-B | | 240-207541 | 240-206908 | 11/19/2015 14:37 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-5-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-5-A | | 240-206833 | 240-206878 | 11/16/2015 13:36 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-5-A | | 240-207234 | 240-206884 | 11/16/2015 08:33 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-5-A | | 240-207234 | 240-206884 | 11/18/2015 10:47 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-5-E | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-5-E | | 240-207364 | 240-206917 | 11/19/2015 01:58 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-5 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

**Lab ID: 240-57930-5 MS** **Client ID: F-VB-UST-P-1**

Sample Date/Time: 11/13/2015 10:00 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:3540C | 240-57930-D-5-C MS | | 240-207541 | 240-206908 | 11/16/2015 09:20 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-5-C MS | | 240-207541 | 240-206908 | 11/19/2015 15:00 | 1 | TAL CAN | JMG |

**Lab ID: 240-57930-5 MSD** **Client ID: F-VB-UST-P-1**

Sample Date/Time: 11/13/2015 10:00 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:3540C | 240-57930-D-5-D MSD | | 240-207541 | 240-206908 | 11/16/2015 09:20 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-5-D MSD | | 240-207541 | 240-206908 | 11/19/2015 15:24 | 1 | TAL CAN | JMG |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Laboratory Chronicle

**Lab ID: 240-57930-6** **Client ID: F-VB-UST-P-2**

Sample Date/Time: 11/13/2015 10:05 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-6-B | | 240-207343 | 240-207393 | 11/18/2015 12:27 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-6-B | | 240-207343 | 240-207393 | 11/18/2015 20:12 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-6-B | | 240-207541 | 240-206887 | 11/16/2015 08:48 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-6-B | | 240-207541 | 240-206887 | 11/19/2015 15:48 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-6-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-6-A | | 240-206833 | 240-206878 | 11/16/2015 14:12 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-6-A | | 240-207234 | 240-206884 | 11/16/2015 08:33 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-6-A | | 240-207234 | 240-206884 | 11/18/2015 11:21 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-6-C | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-6-C | | 240-207364 | 240-206917 | 11/18/2015 22:02 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-6 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

**Lab ID: 240-57930-7** **Client ID: F-VB-UST-P-3**

Sample Date/Time: 11/13/2015 10:10 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-7-B | | 240-207343 | 240-207393 | 11/18/2015 12:27 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-7-B | | 240-207343 | 240-207393 | 11/18/2015 20:33 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-7-B | | 240-207541 | 240-206887 | 11/16/2015 08:48 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-7-B | | 240-207541 | 240-206887 | 11/19/2015 16:12 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-7-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-7-A | | 240-206833 | 240-206878 | 11/16/2015 15:19 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-7-A | | 240-207234 | 240-206884 | 11/16/2015 08:33 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-7-A | | 240-207234 | 240-206884 | 11/18/2015 12:32 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-7-C | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-7-C | | 240-207364 | 240-206917 | 11/18/2015 22:20 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-7 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Laboratory Chronicle

**Lab ID: 240-57930-8** **Client ID: F-VB-UST-P-4**

Sample Date/Time: 11/13/2015 10:15 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-8-A | | 240-207343 | 240-207393 | 11/18/2015 12:27 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-8-A | | 240-207343 | 240-207393 | 11/18/2015 20:55 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-8-A | | 240-207541 | 240-206887 | 11/16/2015 08:48 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-8-A | | 240-207541 | 240-206887 | 11/19/2015 16:36 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-8-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-8-A | | 240-206833 | 240-206878 | 11/16/2015 15:56 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-8-C | | 240-207465 | 240-207121 | 11/17/2015 10:02 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-8-C | | 240-207465 | 240-207121 | 11/19/2015 02:32 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-8-B | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-8-B | | 240-207364 | 240-206917 | 11/18/2015 22:38 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-8 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

**Lab ID: 240-57930-9** **Client ID: F-VB-UST-P-5**

Sample Date/Time: 11/13/2015 10:20 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-9-A | | 240-207343 | 240-207393 | 11/18/2015 12:27 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-9-A | | 240-207343 | 240-207393 | 11/18/2015 21:16 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-9-A | | 240-207541 | 240-206887 | 11/16/2015 08:48 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-9-A | | 240-207541 | 240-206887 | 11/19/2015 17:00 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-9-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-9-A | | 240-206833 | 240-206878 | 11/16/2015 16:30 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-9-C | | 240-207465 | 240-207121 | 11/17/2015 10:02 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-9-C | | 240-207465 | 240-207121 | 11/19/2015 03:04 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-9-B | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-9-B | | 240-207364 | 240-206917 | 11/18/2015 22:56 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-9 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

**Lab ID: 240-57930-9 DU** **Client ID: F-VB-UST-P-5**

Sample Date/Time: 11/13/2015 10:20 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| A:Moisture | 240-57930-A-9 DU | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Laboratory Chronicle

**Lab ID: 240-57930-10** **Client ID: F-VB-UST-P-6**

Sample Date/Time: 11/13/2015 10:25 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-10-A | | 240-207343 | 240-207393 | 11/18/2015 12:27 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-10-A | | 240-207343 | 240-207393 | 11/18/2015 21:38 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-10-A | | 240-207541 | 240-206887 | 11/16/2015 08:48 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-10-A | | 240-207541 | 240-206887 | 11/19/2015 17:23 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-10-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-10-A | | 240-206833 | 240-206878 | 11/16/2015 17:05 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-10-C | | 240-207465 | 240-207121 | 11/17/2015 10:02 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-10-C | | 240-207465 | 240-207121 | 11/19/2015 03:36 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-10-B | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-10-B | | 240-207364 | 240-206917 | 11/18/2015 23:15 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-10 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

**Lab ID: 240-57930-11** **Client ID: F-VS-UST-P-DUP**

Sample Date/Time: 11/13/2015 00:00 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-B-11-A | | 240-207343 | 240-207393 | 11/18/2015 12:30 | 1 | TAL CAN | SAM |
| A:8260C | 240-57930-B-11-A | | 240-207343 | 240-207393 | 11/18/2015 21:59 | 1 | TAL CAN | SAM |
| P:3540C | 240-57930-D-11-A | | 240-207541 | 240-206887 | 11/16/2015 08:48 | 1 | TAL CAN | YDN |
| A:8270D | 240-57930-D-11-A | | 240-207541 | 240-206887 | 11/19/2015 20:10 | 1 | TAL CAN | JMG |
| P:5030A | 240-57930-C-11-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-11-A | | 240-206833 | 240-206878 | 11/16/2015 17:40 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-11-C | | 240-207465 | 240-207121 | 11/17/2015 10:02 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-11-C | | 240-207465 | 240-207121 | 11/19/2015 04:07 | 1 | TAL CAN | DEB |
| P:3540C | 240-57930-D-11-B | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57930-D-11-B | | 240-207364 | 240-206917 | 11/18/2015 23:33 | 1 | TAL CAN | LSH |
| A:Moisture | 240-57930-A-11 | | 240-207016 | | 11/16/2015 15:41 | 1 | TAL CAN | DTN |

**Lab ID: 240-57930-11 MS** **Client ID: F-VS-UST-P-DUP**

Sample Date/Time: 11/13/2015 00:00 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030A | 240-57930-C-11-B MS | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-11-B MS | | 240-206833 | 240-206878 | 11/16/2015 18:14 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-11-D MS | | 240-207465 | 240-207121 | 11/17/2015 10:02 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-11-D MS | | 240-207465 | 240-207121 | 11/19/2015 04:38 | 1 | TAL CAN | DEB |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Laboratory Chronicle

**Lab ID: 240-57930-11 MSD** **Client ID: F-VS-UST-P-DUP**

Sample Date/Time: 11/13/2015 00:00 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030A | 240-57930-C-11-C MSD | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | 240-57930-C-11-C MSD | | 240-206833 | 240-206878 | 11/16/2015 18:49 | 1 | TAL CAN | RTR |
| P:3540C | 240-57930-D-11-E MSD | | 240-207465 | 240-207121 | 11/17/2015 10:02 | 1 | TAL CAN | YDN |
| A:8015C | 240-57930-D-11-E MSD | | 240-207465 | 240-207121 | 11/19/2015 05:10 | 1 | TAL CAN | DEB |

**Lab ID: 240-57930-12** **Client ID: TB-111315**

Sample Date/Time: 11/13/2015 00:00 Received Date/Time: 11/14/2015 09:45

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-57930-A-12 | | 240-207096 | | 11/17/2015 14:51 | 1 | TAL CAN | RJQ |
| A:8260C | 240-57930-A-12 | | 240-207096 | | 11/17/2015 14:51 | 1 | TAL CAN | RJQ |

**Lab ID: MB** **Client ID: N/A**

Sample Date/Time: N/A Received Date/Time: N/A

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | MB 240-207096/6 | | 240-207096 | | 11/17/2015 12:17 | 1 | TAL CAN | RJQ |
| A:8260C | MB 240-207096/6 | | 240-207096 | | 11/17/2015 12:17 | 1 | TAL CAN | RJQ |
| A:8260C | MB 240-207343/6 | | 240-207343 | | 11/18/2015 12:20 | 1 | TAL CAN | SAM |
| A:8260C | MB 240-207560/6 | | 240-207560 | | 11/19/2015 12:29 | 1 | TAL CAN | SAM |
| A:8260C | MB 240-207720/8 | | 240-207720 | | 11/20/2015 05:08 | 1 | TAL CAN | SAM |
| P:3540C | MB 240-206908/21-A | | 240-207287 | 240-206908 | 11/16/2015 09:20 | 1 | TAL CAN | YDN |
| A:8270D | MB 240-206908/21-A | | 240-207287 | 240-206908 | 11/18/2015 08:51 | 1 | TAL CAN | JMG |
| P:3540C | MB 240-206887/23-A | | 240-207541 | 240-206887 | 11/16/2015 08:48 | 1 | TAL CAN | YDN |
| A:8270D | MB 240-206887/23-A | | 240-207541 | 240-206887 | 11/19/2015 11:50 | 1 | TAL CAN | JMG |
| P:5030A | MB 240-206878/1-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | MB 240-206878/1-A | | 240-206833 | 240-206878 | 11/16/2015 10:10 | 1 | TAL CAN | RTR |
| P:5030A | MB 240-207078/1-A | | 240-207037 | 240-207078 | 11/17/2015 07:48 | 1 | TAL CAN | RTR |
| A:8015C | MB 240-207078/1-A | | 240-207037 | 240-207078 | 11/17/2015 09:24 | 1 | TAL CAN | RTR |
| P:5030A | MB 240-207274/1-A | | 240-207246 | 240-207274 | 11/18/2015 06:57 | 1 | TAL CAN | RTR |
| A:8015C | MB 240-207274/1-A | | 240-207246 | 240-207274 | 11/18/2015 15:48 | 1 | TAL CAN | RTR |
| P:3540C | MB 240-206884/23-A | | 240-207234 | 240-206884 | 11/16/2015 08:33 | 1 | TAL CAN | YDN |
| A:8015C | MB 240-206884/23-A | | 240-207234 | 240-206884 | 11/17/2015 22:08 | 1 | TAL CAN | DEB |
| P:3540C | MB 240-207121/7-A | | 240-207465 | 240-207121 | 11/17/2015 10:02 | 1 | TAL CAN | YDN |
| A:8015C | MB 240-207121/7-A | | 240-207465 | 240-207121 | 11/19/2015 01:29 | 1 | TAL CAN | DEB |
| P:3540C | MB 240-206917/21-A | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | MB 240-206917/21-A | | 240-207364 | 240-206917 | 11/19/2015 01:03 | 1 | TAL CAN | LSH |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Laboratory Chronicle

**Lab ID: LCS** **Client ID: N/A**

Sample Date/Time: N/A Received Date/Time: N/A

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | LCS 240-207096/4 | | 240-207096 | | 11/17/2015 11:34 | 1 | TAL CAN | RJQ |
| A:8260C | LCS 240-207096/4 | | 240-207096 | | 11/17/2015 11:34 | 1 | TAL CAN | RJQ |
| A:8260C | LCS 240-207343/5 | | 240-207343 | | 11/18/2015 11:58 | 1 | TAL CAN | SAM |
| A:8260C | LCS 240-207343/7 | | 240-207343 | | 11/18/2015 12:41 | 1 | TAL CAN | SAM |
| A:8260C | LCS 240-207560/5 | | 240-207560 | | 11/19/2015 12:08 | 1 | TAL CAN | SAM |
| A:8260C | LCS 240-207560/7 | | 240-207560 | | 11/19/2015 12:51 | 1 | TAL CAN | SAM |
| A:8260C | LCS 240-207720/6 | | 240-207720 | | 11/20/2015 04:26 | 1 | TAL CAN | SAM |
| P:3540C | LCS 240-206908/22-A | | 240-207287 | 240-206908 | 11/16/2015 09:20 | 1 | TAL CAN | YDN |
| A:8270D | LCS 240-206908/22-A | | 240-207287 | 240-206908 | 11/18/2015 09:15 | 1 | TAL CAN | JMG |
| P:3540C | LCS 240-206887/24-A | | 240-207541 | 240-206887 | 11/16/2015 08:48 | 1 | TAL CAN | YDN |
| A:8270D | LCS 240-206887/24-A | | 240-207541 | 240-206887 | 11/19/2015 12:13 | 1 | TAL CAN | JMG |
| P:5030A | LCS 240-206878/2-A | | 240-206833 | 240-206878 | 11/16/2015 08:12 | 1 | TAL CAN | RTR |
| A:8015C | LCS 240-206878/2-A | | 240-206833 | 240-206878 | 11/16/2015 10:45 | 1 | TAL CAN | RTR |
| P:5030A | LCS 240-207078/2-A | | 240-207037 | 240-207078 | 11/17/2015 07:48 | 1 | TAL CAN | RTR |
| A:8015C | LCS 240-207078/2-A | | 240-207037 | 240-207078 | 11/17/2015 10:01 | 1 | TAL CAN | RTR |
| P:5030A | LCS 240-207274/2-A | | 240-207246 | 240-207274 | 11/18/2015 06:57 | 1 | TAL CAN | RTR |
| A:8015C | LCS 240-207274/2-A | | 240-207246 | 240-207274 | 11/18/2015 16:26 | 1 | TAL CAN | RTR |
| P:3540C | LCS 240-206884/24-A | | 240-207234 | 240-206884 | 11/16/2015 08:33 | 1 | TAL CAN | YDN |
| A:8015C | LCS 240-206884/24-A | | 240-207234 | 240-206884 | 11/17/2015 22:40 | 1 | TAL CAN | DEB |
| P:3540C | LCS 240-207121/8-A | | 240-207465 | 240-207121 | 11/17/2015 10:02 | 1 | TAL CAN | YDN |
| A:8015C | LCS 240-207121/8-A | | 240-207465 | 240-207121 | 11/19/2015 02:01 | 1 | TAL CAN | DEB |
| P:3540C | LCS 240-206917/22-A | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | LCS 240-206917/22-A | | 240-207364 | 240-206917 | 11/19/2015 01:21 | 1 | TAL CAN | LSH |

**Lab ID: MS** **Client ID: N/A**

Sample Date/Time: 11/11/2015 14:30 Received Date/Time: 11/12/2015 14:10

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 200-30727-B-1 MS | | 240-207096 | | 11/17/2015 15:35 | 50 | TAL CAN | RJQ |
| A:8260C | 200-30727-B-1 MS | | 240-207096 | | 11/17/2015 15:35 | 50 | TAL CAN | RJQ |
| P:5030C | 240-57921-A-4-J MS | | 240-207560 | 240-207613 | 11/19/2015 13:00 | 1 | TAL CAN | SAM |
| A:8260C | 240-57921-A-4-J MS | | 240-207560 | 240-207613 | 11/19/2015 20:22 | 1 | TAL CAN | SAM |
| P:5030C | 240-57972-C-3-C MS | | 240-207720 | 240-207709 | 11/20/2015 04:52 | 1 | TAL CAN | TJL2 |
| A:8260C | 240-57972-C-3-C MS | | 240-207720 | 240-207709 | 11/20/2015 08:41 | 1 | TAL CAN | SAM |
| P:5030A | 240-57774-C-8-J MS | | 240-207037 | 240-207078 | 11/17/2015 07:48 | 1 | TAL CAN | RTR |
| A:8015C | 240-57774-C-8-J MS | | 240-207037 | 240-207078 | 11/17/2015 15:35 | 1 | TAL CAN | RTR |
| P:3540C | 240-57880-E-1-B MS | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57880-E-1-B MS | | 240-207364 | 240-206917 | 11/19/2015 02:16 | 1 | TAL CAN | LSH |

Client: Tetra Tech, Inc. Job Number: 240-57930-1

## Laboratory Chronicle

**Lab ID: MSD** **Client ID: N/A**

Sample Date/Time: 11/11/2015 14:30 Received Date/Time: 11/12/2015 14:10

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 200-30727-B-1 MSD | | 240-207096 | | 11/17/2015 15:57 | 50 | TAL CAN | RJQ |
| A:8260C | 200-30727-B-1 MSD | | 240-207096 | | 11/17/2015 15:57 | 50 | TAL CAN | RJQ |
| P:5030C | 240-57921-A-4-K MSD | | 240-207560 | 240-207613 | 11/19/2015 13:00 | 1 | TAL CAN | SAM |
| A:8260C | 240-57921-A-4-K MSD | | 240-207560 | 240-207613 | 11/19/2015 20:43 | 1 | TAL CAN | SAM |
| P:5030C | 240-57972-C-3-D MSD | | 240-207720 | 240-207709 | 11/20/2015 04:52 | 1 | TAL CAN | TJL2 |
| A:8260C | 240-57972-C-3-D MSD | | 240-207720 | 240-207709 | 11/20/2015 09:02 | 1 | TAL CAN | SAM |
| P:5030A | 240-57774-C-8-K MSD | | 240-207037 | 240-207078 | 11/17/2015 07:48 | 1 | TAL CAN | RTR |
| A:8015C | 240-57774-C-8-K MSD | | 240-207037 | 240-207078 | 11/17/2015 16:12 | 1 | TAL CAN | RTR |
| P:3540C | 240-57880-E-1-C MSD | | 240-207364 | 240-206917 | 11/16/2015 09:32 | 1 | TAL CAN | YDN |
| A:8082A | 240-57880-E-1-C MSD | | 240-207364 | 240-206917 | 11/19/2015 02:34 | 1 | TAL CAN | LSH |

**Lab References:**
TAL CAN = TestAmerica Canton

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **ex10PPMSPK_00018** | 04/30/19 | 08/21/15 | MEOH, Lot 147358 | 1000 mL | ex1016/1260st_00001 | 3.8 mL | Aroclor-1016 | 10.0012 ug/mL |
| | | | | | | | Aroclor-1260 | 10.0062 ug/mL |
| | | | | | ex1016/1260st_00002 | 6.2 mL | Aroclor-1016 | 10.0012 ug/mL |
| | | | | | | | Aroclor-1260 | 10.0062 ug/mL |
| .ex1016/1260st_00001 | 04/30/19 | | Restek, Lot A092844 | | (Purchased Reagent) | | Aroclor-1016 | 1000 ug/mL |
| | | | | | | | Aroclor-1260 | 1000 ug/mL |
| .ex1016/1260st_00002 | 05/31/21 | | Restek, Lot A0109056 | | (Purchased Reagent) | | Aroclor-1016 | 1000.2 ug/mL |
| | | | | | | | Aroclor-1260 | 1001 ug/mL |
| **ex2/.2SURRW_00062** | 07/13/17 | 10/26/15 | MEOH, Lot 0000118655 | 2000 mL | EXPESTSURR_00001 | 2 mL | DCB Decachlorobiphenyl | 0.2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.2 ug/mL |
| .EXPESTSURR_00001 | 05/31/21 | | Restek, Lot A0108779 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **exBENZALDEHYD_00031** | 08/31/16 | 10/27/15 | MEOH, Lot 0000099494 | 200 mL | EXLIST1_S11_00001 | 4 mL | Atrazine | 40 ug/mL |
| | | | | | | | Benzaldehyde | 40 ug/mL |
| | | | | | | | Caprolactam | 40 ug/mL |
| .EXLIST1_S11_00001 | 08/31/16 | | Restek, Lot A0108989 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| **exBNASPIKE_00053** | 08/31/16 | 11/02/15 | MEOH, Lot 0000099494 | 200 mL | exLIST1_S1_00004 | 4 mL | 1,1'-Biphenyl | 20 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 20 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 20 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dioxane | 20 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 20 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 40 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 20 ug/mL |
| | | | | | | | 2-Chlorophenol | 20 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2-Methylphenol | 20 ug/mL |
| | | | | | | | 2-Nitroaniline | 20 ug/mL |
| | | | | | | | 2-Nitrophenol | 20 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 20 ug/mL |
| | | | | | | | 3-Methylphenol | 10 ug/mL |
| | | | | | | | 3-Nitroaniline | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 40 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 20 ug/mL |
| | | | | | | | 4-Chloroaniline | 20 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Methylphenol | 10 ug/mL |
| | | | | | | | 4-Nitroaniline | 20 ug/mL |
| | | | | | | | 4-Nitrophenol | 40 ug/mL |
| | | | | | | | Acenaphthene | 20 ug/mL |
| | | | | | | | Acenaphthylene | 20 ug/mL |
| | | | | | | | Acetophenone | 20 ug/mL |
| | | | | | | | Aniline | 20 ug/mL |
| | | | | | | | Anthracene | 20 ug/mL |
| | | | | | | | Azobenzene | 20 ug/mL |
| | | | | | | | Benzo[a]anthracene | 20 ug/mL |
| | | | | | | | Benzo[a]pyrene | 20 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 20 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 20 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 20 ug/mL |
| | | | | | | | Benzyl alcohol | 20 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 20 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 20 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 20 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 20 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 20 ug/mL |
| | | | | | | | Carbazole | 20 ug/mL |
| | | | | | | | Chrysene | 20 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 20 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 20 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 20 ug/mL |
| | | | | | | | Dibenzofuran | 20 ug/mL |
| | | | | | | | Diethyl phthalate | 20 ug/mL |
| | | | | | | | Dimethyl phthalate | 20 ug/mL |
| | | | | | | | Fluoranthene | 20 ug/mL |
| | | | | | | | Fluorene | 20 ug/mL |
| | | | | | | | Hexachlorobenzene | 20 ug/mL |
| | | | | | | | Hexachlorobutadiene | 20 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 20 ug/mL |
| | | | | | | | Hexachloroethane | 20 ug/mL |
| | | | | | | | Hexadecane | 20 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 20 ug/mL |
| | | | | | | | Isophorone | 20 ug/mL |
| | | | | | | | n-Decane | 20 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 40 ug/mL |
| | | | | | | | n-Octadecane | 20 ug/mL |
| | | | | | | | Naphthalene | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Nitrobenzene | 20 ug/mL |
| | | | | | | | Pentachlorophenol | 40 ug/mL |
| | | | | | | | Phenanthrene | 20 ug/mL |
| | | | | | | | Phenol | 20 ug/mL |
| | | | | | | | Pyrene | 20 ug/mL |
| | | | | | | | Pyridine | 20 ug/mL |
| | | | | | EXLIST1_S10_00001 | 4 mL | Benzoic acid | 40 ug/mL |
| | | | | | | | Indene | 40 ug/mL |
| | | | | | exLIST1_S9_00003 | 4 mL | 3,3'-Dichlorobenzidine | 40 ug/mL |
| | | | | | | | Benzidine | 40 ug/mL |
| .exLIST1_S1_00004 | 12/31/16 | | Restek, Lot A0111934 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Methylphenol | 500 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Methylphenol | 500 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| .EXLIST1_S10_00001 | 08/31/16 | | Restek, Lot A0108988 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| .exLIST1_S9_00003 | 01/31/17 | | Restek, Lot A0112567 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **exBNASURR W_00028** | 02/22/18 | 10/15/15 | MEOH, Lot 0000099494 | 2000 mL | SMLIST1 SURR_00003 | 8 mL | 2,4,6-Tribromophenol (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 20 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 20 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 20 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 20 ug/mL |
| .SMLIST1 SURR_00003 | 02/22/18 | Restek, Lot A093638 | | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **exDIESELspkw_00068** | 08/27/17 | 10/05/15 | ACETONE, Lot 0000090987 | 200 mL | exDIESELstd_00003 | 10 mL | Diesel | 2521.45 ug/mL |
| .exDIESELstd_00003 | 04/30/22 | Restek, Lot A0109732 | | | (Purchased Reagent) | | Diesel | 50429 ug/mL |
| **exOTPSURRW_00006** | 12/31/18 | 10/17/15 | MEOH, Lot 0000099494 | 1000 mL | exOTPstd_00002 | 20 mL | o-Terphenyl (Surr) | 200 ug/mL |
| .exOTPstd_00002 | 12/31/18 | Restek, Lot A0110923 | | | (Purchased Reagent) | | o-Terphenyl (Surr) | 10000 ug/mL |
| **Gas Surr Cali_00014** | 11/02/15 | 08/07/15 | MeoH, Lot 1823933 | 10 mL | VGGASURROGATE_00032 | 1 mL | Trifluorotoluene (Surr) | 5 ug/mL |
| .VGGASURROGATE_00032 | 08/22/15 | 05/22/15 | MEOH, Lot 49909 | 50 mL | VGSURRSTKGAS_00019 | 1.25 mL | Trifluorotoluene (Surr) | 50 ug/mL |
| ..VGSURRSTKGAS_00019 | 08/31/16 | ULTRA SCIENTIFIC, Lot CE-1903A | | | (Purchased Reagent) | | Trifluorotoluene (Surr) | 2000 ug/mL |
| **Gas Surr Cali_00015** | 01/12/16 | 10/12/15 | MeoH, Lot 1823933 | 10 mL | VGGASURROGATE_00032 | 1 mL | Trifluorotoluene (Surr) | 5 ug/mL |
| .VGGASURROGATE_00032 | 08/22/15 | 05/22/15 | MEOH, Lot 49909 | 50 mL | VGSURRSTKGAS_00019 | 1.25 mL | Trifluorotoluene (Surr) | 50 ug/mL |
| ..VGSURRSTKGAS_00019 | 08/31/16 | ULTRA SCIENTIFIC, Lot CE-1903A | | | (Purchased Reagent) | | Trifluorotoluene (Surr) | 2000 ug/mL |
| **SG1248@.05ppm_00018** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 0.5 mL | PCB-1248 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.05 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | Restek, Lot A0107141 | | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | Restek, Lot A0108735 | | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.1PPM_00015** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 1 mL | PCB-1248 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.1 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.2ppm_00015** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 2 mL | PCB-1248 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.5ppm_00020** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 5 mL | PCB-1248 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.5 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.5ppm_00020** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 5 mL | Aroclor-1248 | 0.5 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | Aroclor-1248 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | Aroclor-1248 | 1000 ug/mL |
| **SG1248@1.0ppm_00014** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 10 mL | PCB-1248 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@2.0ppm_00014** | 04/22/16 | 08/24/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 20 mL | PCB-1248 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1660@.05PPM_00016** | 04/22/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260_00019 | 0.5 mL | PCB-1016 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.05 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1260 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.05 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.125 mL | DCB Decachlorobiphenyl | 0.0025 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.0025 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | Restek, Lot a109056 | | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | Restek, Lot A0108735 | | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | Restek, Lot A0103185 | | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@0.2ppm_00018** | 04/22/16 | 11/13/15 | HEXANE, Lot 00000101243 | 100 mL | SG1016/1260_00019 | 2 mL | PCB-1016 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.2 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.5 mL | DCB Decachlorobiphenyl | 0.01 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.01 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@0.5PPM_00058** | 04/22/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260_00019 | 5 mL | PCB-1016 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.5 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 1.25 mL | DCB Decachlorobiphenyl | 0.025 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.025 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@0.5PPM_00058** | 04/22/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260_00019 | 5 mL | Aroclor-1016 | 0.5 ug/mL |
| | | | | | | | Aroclor-1260 | 0.5 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | Aroclor-1016 | 10 ug/mL |
| | | | | | | | Aroclor-1260 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | Aroclor-1016 | 1000 ug/mL |
| | | | | | | | Aroclor-1260 | 1000 ug/mL |
| **SG1660@1.0PPM_00016** | 04/22/16 | 11/13/15 | HEXANE, Lot 00000101243 | 100 mL | SG1016/1260_00019 | 10 mL | PCB-1016 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 2.5 mL | DCB Decachlorobiphenyl | 0.05 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.05 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@2.0PPM_00015** | 04/22/16 | 11/13/15 | HEXANE, Lot 00000101243 | 100 mL | SG1016/1260_00019 | 20 mL | PCB-1016 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 5 mL | DCB Decachlorobiphenyl | 0.1 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.1 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660ICV@1.0_00015** | 04/21/16 | 11/03/15 | HEXANE, Lot 0000101243 | 10 mL | SGPCBIS STOCk_00003 | 0.05 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1660ICV@1.0_00015** | 04/21/16 | 11/03/15 | HEXANE, Lot 0000101243 | 10 mL | SG1660ICV@100_00005 | 0.1 mL | Aroclor-1016 | 1 ug/mL |
| | | | | | | | Aroclor-1260 | 1 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.25 mL | DCB Decachlorobiphenyl | 0.05 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.05 ug/mL |
| .SG1660ICV@100_00005 | 10/28/16 | 10/28/15 | HEXANE, Lot 0000101243 | 10 mL | SG 1660 ICV_00012 | 1 mL | Aroclor-1016 | 100 ug/mL |
| | | | | | | | Aroclor-1260 | 100 ug/mL |
| ..SG 1660 ICV_00012 | 01/30/22 | | restek, Lot a0114674 | | (Purchased Reagent) | | Aroclor-1016 | 1000 ug/mL |
| | | | | | | | Aroclor-1260 | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660STD@0.1_00016** | 04/22/16 | 11/13/15 | HEXANE, Lot 154006 | 100 mL | SG1016/1260_00019 | 1 mL | PCB-1016 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.1 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.25 mL | DCB Decachlorobiphenyl | 0.005 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.005 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG2154@0.05PP_00010** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 0.5 mL | PCB-1221 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.05 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@0.2PPM_00012** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 2 mL | PCB-1221 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.2 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1254 Peak 5 | 0.2 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@0.5PPM_00021** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 5 mL | PCB-1221 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.5 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **SG2154@0.5PPM_00021** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 5 mL | Aroclor-1221 | 0.5 ug/mL |
| | | | | | | | Aroclor-1254 | 0.5 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | Aroclor-1221 | 10 ug/mL |
| | | | | | | | Aroclor-1254 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | Aroclor-1221 | 1000 ug/mL |
| | | | | | | | Aroclor-1254 | 1000 ug/mL |
| **SG2154@1.0PPM_00014** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 10 mL | PCB-1221 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@2.0PPM_00011** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 20 mL | PCB-1221 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 2 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@O.1PPM_00013** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 1 mL | PCB-1221 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.1 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@.05PPM_00004** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00003 | 0.5 mL | PCB-1232 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 50 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1262 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 50 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00003 | 09/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/62 STK_00001 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SG32/62 STK_00001 | 04/30/21 | Restelk, Lot A0108264 | | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | Restek, Lot A0108735 | | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@.2PPM_00005** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 2 mL | PCB-1232 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 200 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@0.1PPM_00006** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 1 mL | PCB-1232 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 100 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@0.5PPM_00010** | 04/22/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00003 | 5 mL | PCB-1232 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 500 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1232 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 500 ug/L |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00003 | 09/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/62 STK_00001 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SG32/62 STK_00001 | 04/30/21 | | Restelk, Lot A0108264 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@0.5PPM_00010** | 04/22/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00003 | 5 mL | Aroclor-1232 | 500 ug/L |
| | | | | | | | PCB-1262 | 500 ug/L |
| .SG32/6210PPM_00003 | 09/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/62 STK_00001 | 1 mL | Aroclor-1232 | 10000 ug/L |
| | | | | | | | PCB-1262 | 10000 ug/L |
| ..SG32/62 STK_00001 | 04/30/21 | | Restelk, Lot A0108264 | | (Purchased Reagent) | | Aroclor-1232 | 1000 ug/mL |
| | | | | | | | PCB-1262 | 1000 ug/mL |
| **SG3262@1.0PPM_00005** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 10 mL | PCB-1232 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@2.0ppm_00004** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 20 mL | PCB-1232 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 2000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG42/68@1.0PP_00008** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 10 mL | PCB-1242 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@.05PPM_00005** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 0.5 mL | PCB-1242 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 50 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1268 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 50 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@.1PPM_00006** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 1 mL | PCB-1242 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 100 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@.2PPM_00005** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 2 mL | PCB-1242 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 200 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **SG4268@0.5PPM_00007** | 12/24/15 | 06/24/15 | HEXANE, Lot 0000086560 | 100 mL | SG42/6810PPM_00002 | 5 mL | PCB-1242 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 500 ug/L |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | Restek, Lot A090182 | | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | Restek, Lot A097795 | | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | Restek, Lot A0110081 | | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@0.5PPM_00011** | 04/07/16 | 10/07/15 | HEXANE, Lot 154006 | 100 mL | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | Restek, Lot A0110081 | | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@0.5PPM_00011** | 04/07/16 | 10/07/15 | HEXANE, Lot 154006 | 100 mL | SG42/6810PPM_00003 | 5 mL | Aroclor-1242 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 500 ug/L |
| | | | | | | | PCB-1268 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 500 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1268 Peak 5 | 500 ug/L |
| .SG42/6810PPM_00003 | 10/07/16 | 10/07/15 | HEXANE, Lot 154006 | 100 mL | SG 42/68 STK_00001 | 1 mL | Aroclor-1242 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | | | PCB-1268 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG 42/68 STK_00001 | 04/30/21 | | RESTEK, Lot A0108269 | | (Purchased Reagent) | | Aroclor-1242 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1268 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| **SG4268@2.0PPM_00004** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 20 mL | PCB-1242 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 2000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SGDROCAL_00020** | 10/05/16 | 10/05/15 | MECL2, Lot EXDCM_00187 | 25 mL | SGDROCALSTK_00019 | 5 mL | Diesel | 10002.5 ug/mL |
| | | | | | SGDROSURROTP_00025 | 1 mL | o-Terphenyl (Surr) | 800 ug/mL |
| | | | | | SGDROSURROTP_00026 | 1 mL | o-Terphenyl (Surr) | 800 ug/mL |
| .SGDROCALSTK_00019 | 04/30/22 | | Restek, Lot A0109723 | | (Purchased Reagent) | | Diesel | 50005.8 ug/mL |
| .SGDROSURROTP_00025 | 09/30/18 | | Restek, Lot A0109488 | | (Purchased Reagent) | | o-Terphenyl (Surr) | 10062.7 ug/mL |
| .SGDROSURROTP_00026 | 09/30/18 | | Restek, Lot A0109488 | | (Purchased Reagent) | | o-Terphenyl (Surr) | 10062.7 ug/mL |
| **SGDROICV_00015** | 08/26/16 | 08/26/15 | MECL2, Lot 2091477 | 25 mL | SGDROICVSTK_00008 | 1 mL | Diesel | 2006 ug/mL |
| .SGDROICVSTK_00008 | 08/31/18 | | Ultra, Lot CL-2929 | | (Purchased Reagent) | | Diesel | 50150 ug/mL |
| **SGPCBISTD_00003** | 04/22/16 | 10/13/15 | HEXANE, Lot 154006 | 100 mL | SGPCBIS STOCk_00003 | 10 mL | 1-Bromo-2-nitrobenzene | 1 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SMIS80PPMW_00013** | 07/10/16 | 07/10/15 | MECL2, Lot 0000103549 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| .SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| **SMLIST1 L1 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1PAH ST_00004 | 2 uL | 1,1'-Biphenyl | 0.1 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 0.1 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 0.1 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 0.1 ug/mL |
| | | | | | | | Acenaphthene | 0.1 ug/mL |
| | | | | | | | Acenaphthylene | 0.1 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Anthracene | 0.1 ug/mL |
| | | | | | | | Benzo[a]anthracene | 0.1 ug/mL |
| | | | | | | | Benzo[a]pyrene | 0.1 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 0.1 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 0.1 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 0.1 ug/mL |
| | | | | | | | Chrysene | 0.1 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 0.1 ug/mL |
| | | | | | | | Dibenzofuran | 0.1 ug/mL |
| | | | | | | | Fluoranthene | 0.1 ug/mL |
| | | | | | | | Fluorene | 0.1 ug/mL |
| | | | | | | | Hexachlorobenzene | 0.1 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 0.1 ug/mL |
| | | | | | | | Naphthalene | 0.1 ug/mL |
| | | | | | | | Phenanthrene | 0.1 ug/mL |
| | | | | | | | Pyrene | 0.1 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 0.1 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 0.1 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 0.1 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1PAH ST_00004 | 05/31/16 | 01/26/15 | MECL2, Lot 0000093871 | 3 mL | SMLIST1 PAH_00004 | 300 uL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 60 uL | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 PAH_00004 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L2 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 10 uL | 1,1'-Biphenyl | 0.5 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 0.5 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 0.5 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 0.5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,4-Dioxane | 0.5 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 0.5 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 0.5 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 1 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 0.5 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 0.5 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 0.5 ug/mL |
| | | | | | | | 2-Chlorophenol | 0.5 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 0.5 ug/mL |
| | | | | | | | 2-Methylphenol | 0.5 ug/mL |
| | | | | | | | 2-Nitroaniline | 0.5 ug/mL |
| | | | | | | | 2-Nitrophenol | 0.5 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 0.5 ug/mL |
| | | | | | | | 3-Nitroaniline | 0.5 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 1 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 0.5 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 0.5 ug/mL |
| | | | | | | | 4-Chloroaniline | 0.5 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 0.5 ug/mL |
| | | | | | | | 4-Nitroaniline | 0.5 ug/mL |
| | | | | | | | 4-Nitrophenol | 1 ug/mL |
| | | | | | | | Acenaphthene | 0.5 ug/mL |
| | | | | | | | Acenaphthylene | 0.5 ug/mL |
| | | | | | | | Acetophenone | 0.5 ug/mL |
| | | | | | | | Aniline | 0.5 ug/mL |
| | | | | | | | Anthracene | 0.5 ug/mL |
| | | | | | | | Azobenzene | 0.5 ug/mL |
| | | | | | | | Benzo[a]anthracene | 0.5 ug/mL |
| | | | | | | | Benzo[a]pyrene | 0.5 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 0.5 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 0.5 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 0.5 ug/mL |
| | | | | | | | Benzyl alcohol | 0.5 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 0.5 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 0.5 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 0.5 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 0.5 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 0.5 ug/mL |
| | | | | | | | Carbazole | 0.5 ug/mL |
| | | | | | | | Chrysene | 0.5 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 0.5 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 0.5 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 0.5 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Dibenzofuran | 0.5 ug/mL |
| | | | | | | | Diethyl phthalate | 0.5 ug/mL |
| | | | | | | | Dimethyl phthalate | 0.5 ug/mL |
| | | | | | | | Fluoranthene | 0.5 ug/mL |
| | | | | | | | Fluorene | 0.5 ug/mL |
| | | | | | | | Hexachlorobenzene | 0.5 ug/mL |
| | | | | | | | Hexachlorobutadiene | 0.5 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 0.5 ug/mL |
| | | | | | | | Hexachloroethane | 0.5 ug/mL |
| | | | | | | | Hexadecane | 0.5 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 0.5 ug/mL |
| | | | | | | | Isophorone | 0.5 ug/mL |
| | | | | | | | n-Decane | 0.5 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 0.5 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 0.5 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 1 ug/mL |
| | | | | | | | n-Octadecane | 0.5 ug/mL |
| | | | | | | | Naphthalene | 0.5 ug/mL |
| | | | | | | | Nitrobenzene | 0.5 ug/mL |
| | | | | | | | Pentachlorophenol | 1 ug/mL |
| | | | | | | | Phenanthrene | 0.5 ug/mL |
| | | | | | | | Phenol | 0.5 ug/mL |
| | | | | | | | Pyrene | 0.5 ug/mL |
| | | | | | | | Pyridine | 0.5 ug/mL |
| | | | | | | | Benzoic acid | 1 ug/mL |
| | | | | | | | Indene | 1 ug/mL |
| | | | | | | | Atrazine | 1 ug/mL |
| | | | | | | | Benzaldehyde | 1 ug/mL |
| | | | | | | | Caprolactam | 1 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 1 ug/mL |
| | | | | | | | Benzidine | 1 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 0.5 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 0.5 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 0.5 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 0.5 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 0.5 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 0.5 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L3 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 20 uL | 1,1'-Biphenyl | 1 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1 ug/mL |
| | | | | | | | 1,4-Dioxane | 1 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1 ug/mL |
| | | | | | | | 2-Chlorophenol | 1 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1 ug/mL |
| | | | | | | | 2-Methylphenol | 1 ug/mL |
| | | | | | | | 2-Nitroaniline | 1 ug/mL |
| | | | | | | | 2-Nitrophenol | 1 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1 ug/mL |
| | | | | | | | 3-Nitroaniline | 1 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1 ug/mL |
| | | | | | | | 4-Chloroaniline | 1 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1 ug/mL |
| | | | | | | | 4-Nitroaniline | 1 ug/mL |
| | | | | | | | 4-Nitrophenol | 2 ug/mL |
| | | | | | | | Acenaphthene | 1 ug/mL |
| | | | | | | | Acenaphthylene | 1 ug/mL |
| | | | | | | | Acetophenone | 1 ug/mL |
| | | | | | | | Aniline | 1 ug/mL |
| | | | | | | | Anthracene | 1 ug/mL |
| | | | | | | | Azobenzene | 1 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1 ug/mL |
| | | | | | | | Benzyl alcohol | 1 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | bis (2-chloroisopropyl) ether | 1 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1 ug/mL |
| | | | | | | | Carbazole | 1 ug/mL |
| | | | | | | | Chrysene | 1 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1 ug/mL |
| | | | | | | | Dibenzofuran | 1 ug/mL |
| | | | | | | | Diethyl phthalate | 1 ug/mL |
| | | | | | | | Dimethyl phthalate | 1 ug/mL |
| | | | | | | | Fluoranthene | 1 ug/mL |
| | | | | | | | Fluorene | 1 ug/mL |
| | | | | | | | Hexachlorobenzene | 1 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1 ug/mL |
| | | | | | | | Hexachloroethane | 1 ug/mL |
| | | | | | | | Hexadecane | 1 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1 ug/mL |
| | | | | | | | Isophorone | 1 ug/mL |
| | | | | | | | n-Decane | 1 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2 ug/mL |
| | | | | | | | n-Octadecane | 1 ug/mL |
| | | | | | | | Naphthalene | 1 ug/mL |
| | | | | | | | Nitrobenzene | 1 ug/mL |
| | | | | | | | Pentachlorophenol | 2 ug/mL |
| | | | | | | | Phenanthrene | 1 ug/mL |
| | | | | | | | Phenol | 1 ug/mL |
| | | | | | | | Pyrene | 1 ug/mL |
| | | | | | | | Pyridine | 1 ug/mL |
| | | | | | | | Benzoic acid | 2 ug/mL |
| | | | | | | | Indene | 2 ug/mL |
| | | | | | | | Atrazine | 2 ug/mL |
| | | | | | | | Benzaldehyde | 2 ug/mL |
| | | | | | | | Caprolactam | 2 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 2 ug/mL |
| | | | | | | | Benzidine | 2 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 1 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 1 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 1 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 1 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 1 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 1 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L4 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 40 uL | 1,1'-Biphenyl | 2 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 2 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 2 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 2 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2 ug/mL |
| | | | | | | | 1,4-Dioxane | 2 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 2 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 2 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 2 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 2 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 2 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 2 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 4 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 2 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 2 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 2 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 2 ug/mL |
| | | | | | | | 2-Chlorophenol | 2 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 2 ug/mL |
| | | | | | | | 2-Methylphenol | 2 ug/mL |
| | | | | | | | 2-Nitroaniline | 2 ug/mL |
| | | | | | | | 2-Nitrophenol | 2 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 2 ug/mL |
| | | | | | | | 3-Nitroaniline | 2 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 4 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 2 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 2 ug/mL |
| | | | | | | | 4-Chloroaniline | 2 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 2 ug/mL |
| | | | | | | | 4-Nitroaniline | 2 ug/mL |
| | | | | | | | 4-Nitrophenol | 4 ug/mL |
| | | | | | | | Acenaphthene | 2 ug/mL |
| | | | | | | | Acenaphthylene | 2 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acetophenone | 2 ug/mL |
| | | | | | | | Aniline | 2 ug/mL |
| | | | | | | | Anthracene | 2 ug/mL |
| | | | | | | | Azobenzene | 2 ug/mL |
| | | | | | | | Benzo[a]anthracene | 2 ug/mL |
| | | | | | | | Benzo[a]pyrene | 2 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 2 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 2 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 2 ug/mL |
| | | | | | | | Benzyl alcohol | 2 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 2 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 2 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 2 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 2 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 2 ug/mL |
| | | | | | | | Carbazole | 2 ug/mL |
| | | | | | | | Chrysene | 2 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 2 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 2 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 2 ug/mL |
| | | | | | | | Dibenzofuran | 2 ug/mL |
| | | | | | | | Diethyl phthalate | 2 ug/mL |
| | | | | | | | Dimethyl phthalate | 2 ug/mL |
| | | | | | | | Fluoranthene | 2 ug/mL |
| | | | | | | | Fluorene | 2 ug/mL |
| | | | | | | | Hexachlorobenzene | 2 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 2 ug/mL |
| | | | | | | | Hexachloroethane | 2 ug/mL |
| | | | | | | | Hexadecane | 2 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 2 ug/mL |
| | | | | | | | Isophorone | 2 ug/mL |
| | | | | | | | n-Decane | 2 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 2 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 2 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 4 ug/mL |
| | | | | | | | n-Octadecane | 2 ug/mL |
| | | | | | | | Naphthalene | 2 ug/mL |
| | | | | | | | Nitrobenzene | 2 ug/mL |
| | | | | | | | Pentachlorophenol | 4 ug/mL |
| | | | | | | | Phenanthrene | 2 ug/mL |
| | | | | | | | Phenol | 2 ug/mL |
| | | | | | | | Pyrene | 2 ug/mL |
| | | | | | | | Pyridine | 2 ug/mL |
| | | | | | | | Benzoic acid | 4 ug/mL |
| | | | | | | | Indene | 4 ug/mL |
| | | | | | | | Atrazine | 4 ug/mL |
| | | | | | | | Benzaldehyde | 4 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Caprolactam | 4 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 4 ug/mL |
| | | | | | | | Benzidine | 4 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 2 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 2 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 2 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 2 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 2 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 2 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L5 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 100 uL | 1,1'-Biphenyl | 5 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 5 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 5 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 5 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 5 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 5 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 5 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 5 ug/mL |
| | | | | | | | 1,4-Dioxane | 5 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 5 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 5 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 5 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 5 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 5 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 5 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 10 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 5 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 5 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 5 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 5 ug/mL |
| | | | | | | | 2-Chlorophenol | 5 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 5 ug/mL |
| | | | | | | | 2-Methylphenol | 5 ug/mL |
| | | | | | | | 2-Nitroaniline | 5 ug/mL |
| | | | | | | | 2-Nitrophenol | 5 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 3-Nitroaniline | 5 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 10 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 5 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 5 ug/mL |
| | | | | | | | 4-Chloroaniline | 5 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 5 ug/mL |
| | | | | | | | 4-Nitroaniline | 5 ug/mL |
| | | | | | | | 4-Nitrophenol | 10 ug/mL |
| | | | | | | | Acenaphthene | 5 ug/mL |
| | | | | | | | Acenaphthylene | 5 ug/mL |
| | | | | | | | Acetophenone | 5 ug/mL |
| | | | | | | | Aniline | 5 ug/mL |
| | | | | | | | Anthracene | 5 ug/mL |
| | | | | | | | Azobenzene | 5 ug/mL |
| | | | | | | | Benzo[a]anthracene | 5 ug/mL |
| | | | | | | | Benzo[a]pyrene | 5 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 5 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 5 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 5 ug/mL |
| | | | | | | | Benzyl alcohol | 5 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 5 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 5 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 5 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 5 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 5 ug/mL |
| | | | | | | | Carbazole | 5 ug/mL |
| | | | | | | | Chrysene | 5 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 5 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 5 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 5 ug/mL |
| | | | | | | | Dibenzofuran | 5 ug/mL |
| | | | | | | | Diethyl phthalate | 5 ug/mL |
| | | | | | | | Dimethyl phthalate | 5 ug/mL |
| | | | | | | | Fluoranthene | 5 ug/mL |
| | | | | | | | Fluorene | 5 ug/mL |
| | | | | | | | Hexachlorobenzene | 5 ug/mL |
| | | | | | | | Hexachlorobutadiene | 5 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 5 ug/mL |
| | | | | | | | Hexachloroethane | 5 ug/mL |
| | | | | | | | Hexadecane | 5 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 5 ug/mL |
| | | | | | | | Isophorone | 5 ug/mL |
| | | | | | | | n-Decane | 5 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 5 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 5 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 10 ug/mL |
| | | | | | | | n-Octadecane | 5 ug/mL |
| | | | | | | | Naphthalene | 5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Nitrobenzene | 5 ug/mL |
| | | | | | | | Pentachlorophenol | 10 ug/mL |
| | | | | | | | Phenanthrene | 5 ug/mL |
| | | | | | | | Phenol | 5 ug/mL |
| | | | | | | | Pyrene | 5 ug/mL |
| | | | | | | | Pyridine | 5 ug/mL |
| | | | | | | | Benzoic acid | 10 ug/mL |
| | | | | | | | Indene | 10 ug/mL |
| | | | | | | | Atrazine | 10 ug/mL |
| | | | | | | | Benzaldehyde | 10 ug/mL |
| | | | | | | | Caprolactam | 10 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 10 ug/mL |
| | | | | | | | Benzidine | 10 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 5 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L6 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 200 uL | 1,1'-Biphenyl | 10 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 10 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 10 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 10 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 10 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 10 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 10 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 10 ug/mL |
| | | | | | | | 1,4-Dioxane | 10 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 10 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 10 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 10 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,4-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 10 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 10 ug/mL |
| | | | | | | | 2-Chlorophenol | 10 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 10 ug/mL |
| | | | | | | | 2-Methylphenol | 10 ug/mL |
| | | | | | | | 2-Nitroaniline | 10 ug/mL |
| | | | | | | | 2-Nitrophenol | 10 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 10 ug/mL |
| | | | | | | | 3-Nitroaniline | 10 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 20 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 10 ug/mL |
| | | | | | | | 4-Chloroaniline | 10 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Nitroaniline | 10 ug/mL |
| | | | | | | | 4-Nitrophenol | 20 ug/mL |
| | | | | | | | Acenaphthene | 10 ug/mL |
| | | | | | | | Acenaphthylene | 10 ug/mL |
| | | | | | | | Acetophenone | 10 ug/mL |
| | | | | | | | Aniline | 10 ug/mL |
| | | | | | | | Anthracene | 10 ug/mL |
| | | | | | | | Azobenzene | 10 ug/mL |
| | | | | | | | Benzo[a]anthracene | 10 ug/mL |
| | | | | | | | Benzo[a]pyrene | 10 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 10 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 10 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 10 ug/mL |
| | | | | | | | Benzyl alcohol | 10 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 10 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 10 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 10 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 10 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 10 ug/mL |
| | | | | | | | Carbazole | 10 ug/mL |
| | | | | | | | Chrysene | 10 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 10 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 10 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 10 ug/mL |
| | | | | | | | Dibenzofuran | 10 ug/mL |
| | | | | | | | Diethyl phthalate | 10 ug/mL |
| | | | | | | | Dimethyl phthalate | 10 ug/mL |
| | | | | | | | Fluoranthene | 10 ug/mL |
| | | | | | | | Fluorene | 10 ug/mL |
| | | | | | | | Hexachlorobenzene | 10 ug/mL |
| | | | | | | | Hexachlorobutadiene | 10 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Hexachloroethane | 10 ug/mL |
| | | | | | | | Hexadecane | 10 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 10 ug/mL |
| | | | | | | | Isophorone | 10 ug/mL |
| | | | | | | | n-Decane | 10 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 10 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 10 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 20 ug/mL |
| | | | | | | | n-Octadecane | 10 ug/mL |
| | | | | | | | Naphthalene | 10 ug/mL |
| | | | | | | | Nitrobenzene | 10 ug/mL |
| | | | | | | | Pentachlorophenol | 20 ug/mL |
| | | | | | | | Phenanthrene | 10 ug/mL |
| | | | | | | | Phenol | 10 ug/mL |
| | | | | | | | Pyrene | 10 ug/mL |
| | | | | | | | Pyridine | 10 ug/mL |
| | | | | | | | Benzoic acid | 20 ug/mL |
| | | | | | | | Indene | 20 ug/mL |
| | | | | | | | Atrazine | 20 ug/mL |
| | | | | | | | Benzaldehyde | 20 ug/mL |
| | | | | | | | Caprolactam | 20 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 20 ug/mL |
| | | | | | | | Benzidine | 20 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 10 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 10 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 10 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 10 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L7 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 300 uL | 1,1'-Biphenyl | 15 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 15 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 15 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 15 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 15 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 15 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,3-Dinitrobenzene | 15 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 15 ug/mL |
| | | | | | | | 1,4-Dioxane | 15 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 15 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 15 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 15 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 15 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 15 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 15 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 30 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 15 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 15 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 15 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 15 ug/mL |
| | | | | | | | 2-Chlorophenol | 15 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 15 ug/mL |
| | | | | | | | 2-Methylphenol | 15 ug/mL |
| | | | | | | | 2-Nitroaniline | 15 ug/mL |
| | | | | | | | 2-Nitrophenol | 15 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 15 ug/mL |
| | | | | | | | 3-Nitroaniline | 15 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 30 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 15 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 15 ug/mL |
| | | | | | | | 4-Chloroaniline | 15 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 15 ug/mL |
| | | | | | | | 4-Nitroaniline | 15 ug/mL |
| | | | | | | | 4-Nitrophenol | 30 ug/mL |
| | | | | | | | Acenaphthene | 15 ug/mL |
| | | | | | | | Acenaphthylene | 15 ug/mL |
| | | | | | | | Acetophenone | 15 ug/mL |
| | | | | | | | Aniline | 15 ug/mL |
| | | | | | | | Anthracene | 15 ug/mL |
| | | | | | | | Azobenzene | 15 ug/mL |
| | | | | | | | Benzo[a]anthracene | 15 ug/mL |
| | | | | | | | Benzo[a]pyrene | 15 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 15 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 15 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 15 ug/mL |
| | | | | | | | Benzyl alcohol | 15 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 15 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 15 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 15 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 15 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 15 ug/mL |
| | | | | | | | Carbazole | 15 ug/mL |
| | | | | | | | Chrysene | 15 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 15 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Di-n-octyl phthalate | 15 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 15 ug/mL |
| | | | | | | | Dibenzofuran | 15 ug/mL |
| | | | | | | | Diethyl phthalate | 15 ug/mL |
| | | | | | | | Dimethyl phthalate | 15 ug/mL |
| | | | | | | | Fluoranthene | 15 ug/mL |
| | | | | | | | Fluorene | 15 ug/mL |
| | | | | | | | Hexachlorobenzene | 15 ug/mL |
| | | | | | | | Hexachlorobutadiene | 15 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 15 ug/mL |
| | | | | | | | Hexachloroethane | 15 ug/mL |
| | | | | | | | Hexadecane | 15 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 15 ug/mL |
| | | | | | | | Isophorone | 15 ug/mL |
| | | | | | | | n-Decane | 15 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 15 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 15 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 30 ug/mL |
| | | | | | | | n-Octadecane | 15 ug/mL |
| | | | | | | | Naphthalene | 15 ug/mL |
| | | | | | | | Nitrobenzene | 15 ug/mL |
| | | | | | | | Pentachlorophenol | 30 ug/mL |
| | | | | | | | Phenanthrene | 15 ug/mL |
| | | | | | | | Phenol | 15 ug/mL |
| | | | | | | | Pyrene | 15 ug/mL |
| | | | | | | | Pyridine | 15 ug/mL |
| | | | | | | | Benzoic acid | 30 ug/mL |
| | | | | | | | Indene | 30 ug/mL |
| | | | | | | | Atrazine | 30 ug/mL |
| | | | | | | | Benzaldehyde | 30 ug/mL |
| | | | | | | | Caprolactam | 30 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 30 ug/mL |
| | | | | | | | Benzidine | 30 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 15 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 15 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 15 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 15 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 15 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 15 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L8 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 400 uL | 1,1'-Biphenyl | 20 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 20 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 20 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dioxane | 20 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 20 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 40 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 20 ug/mL |
| | | | | | | | 2-Chlorophenol | 20 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2-Methylphenol | 20 ug/mL |
| | | | | | | | 2-Nitroaniline | 20 ug/mL |
| | | | | | | | 2-Nitrophenol | 20 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 20 ug/mL |
| | | | | | | | 3-Nitroaniline | 20 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 40 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 20 ug/mL |
| | | | | | | | 4-Chloroaniline | 20 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Nitroaniline | 20 ug/mL |
| | | | | | | | 4-Nitrophenol | 40 ug/mL |
| | | | | | | | Acenaphthene | 20 ug/mL |
| | | | | | | | Acenaphthylene | 20 ug/mL |
| | | | | | | | Acetophenone | 20 ug/mL |
| | | | | | | | Aniline | 20 ug/mL |
| | | | | | | | Anthracene | 20 ug/mL |
| | | | | | | | Azobenzene | 20 ug/mL |
| | | | | | | | Benzo[a]anthracene | 20 ug/mL |
| | | | | | | | Benzo[a]pyrene | 20 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 20 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 20 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[k]fluoranthene | 20 ug/mL |
| | | | | | | | Benzyl alcohol | 20 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 20 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 20 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 20 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 20 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 20 ug/mL |
| | | | | | | | Carbazole | 20 ug/mL |
| | | | | | | | Chrysene | 20 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 20 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 20 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 20 ug/mL |
| | | | | | | | Dibenzofuran | 20 ug/mL |
| | | | | | | | Diethyl phthalate | 20 ug/mL |
| | | | | | | | Dimethyl phthalate | 20 ug/mL |
| | | | | | | | Fluoranthene | 20 ug/mL |
| | | | | | | | Fluorene | 20 ug/mL |
| | | | | | | | Hexachlorobenzene | 20 ug/mL |
| | | | | | | | Hexachlorobutadiene | 20 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 20 ug/mL |
| | | | | | | | Hexachloroethane | 20 ug/mL |
| | | | | | | | Hexadecane | 20 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 20 ug/mL |
| | | | | | | | Isophorone | 20 ug/mL |
| | | | | | | | n-Decane | 20 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 40 ug/mL |
| | | | | | | | n-Octadecane | 20 ug/mL |
| | | | | | | | Naphthalene | 20 ug/mL |
| | | | | | | | Nitrobenzene | 20 ug/mL |
| | | | | | | | Pentachlorophenol | 40 ug/mL |
| | | | | | | | Phenanthrene | 20 ug/mL |
| | | | | | | | Phenol | 20 ug/mL |
| | | | | | | | Pyrene | 20 ug/mL |
| | | | | | | | Pyridine | 20 ug/mL |
| | | | | | | | Benzoic acid | 40 ug/mL |
| | | | | | | | Indene | 40 ug/mL |
| | | | | | | | Atrazine | 40 ug/mL |
| | | | | | | | Benzaldehyde | 40 ug/mL |
| | | | | | | | Caprolactam | 40 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 40 ug/mL |
| | | | | | | | Benzidine | 40 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 20 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 20 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Terphenyl-d14 (Surr) | 20 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L9 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 500 uL | 1,1'-Biphenyl | 25 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 25 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 25 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 25 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 25 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 25 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 25 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 25 ug/mL |
| | | | | | | | 1,4-Dioxane | 25 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 25 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 25 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 25 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 25 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 25 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 25 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 50 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 25 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 25 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 25 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 25 ug/mL |
| | | | | | | | 2-Chlorophenol | 25 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 25 ug/mL |
| | | | | | | | 2-Methylphenol | 25 ug/mL |
| | | | | | | | 2-Nitroaniline | 25 ug/mL |
| | | | | | | | 2-Nitrophenol | 25 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 25 ug/mL |
| | | | | | | | 3-Nitroaniline | 25 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 50 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 25 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 25 ug/mL |
| | | | | | | | 4-Chloroaniline | 25 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 25 ug/mL |
| | | | | | | | 4-Nitroaniline | 25 ug/mL |
| | | | | | | | 4-Nitrophenol | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acenaphthene | 25 ug/mL |
| | | | | | | | Acenaphthylene | 25 ug/mL |
| | | | | | | | Acetophenone | 25 ug/mL |
| | | | | | | | Aniline | 25 ug/mL |
| | | | | | | | Anthracene | 25 ug/mL |
| | | | | | | | Azobenzene | 25 ug/mL |
| | | | | | | | Benzo[a]anthracene | 25 ug/mL |
| | | | | | | | Benzo[a]pyrene | 25 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 25 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 25 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 25 ug/mL |
| | | | | | | | Benzyl alcohol | 25 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 25 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 25 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 25 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 25 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 25 ug/mL |
| | | | | | | | Carbazole | 25 ug/mL |
| | | | | | | | Chrysene | 25 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 25 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 25 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 25 ug/mL |
| | | | | | | | Dibenzofuran | 25 ug/mL |
| | | | | | | | Diethyl phthalate | 25 ug/mL |
| | | | | | | | Dimethyl phthalate | 25 ug/mL |
| | | | | | | | Fluoranthene | 25 ug/mL |
| | | | | | | | Fluorene | 25 ug/mL |
| | | | | | | | Hexachlorobenzene | 25 ug/mL |
| | | | | | | | Hexachlorobutadiene | 25 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 25 ug/mL |
| | | | | | | | Hexachloroethane | 25 ug/mL |
| | | | | | | | Hexadecane | 25 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 25 ug/mL |
| | | | | | | | Isophorone | 25 ug/mL |
| | | | | | | | n-Decane | 25 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 25 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 25 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 50 ug/mL |
| | | | | | | | n-Octadecane | 25 ug/mL |
| | | | | | | | Naphthalene | 25 ug/mL |
| | | | | | | | Nitrobenzene | 25 ug/mL |
| | | | | | | | Pentachlorophenol | 50 ug/mL |
| | | | | | | | Phenanthrene | 25 ug/mL |
| | | | | | | | Phenol | 25 ug/mL |
| | | | | | | | Pyrene | 25 ug/mL |
| | | | | | | | Pyridine | 25 ug/mL |
| | | | | | | | Benzoic acid | 50 ug/mL |
| | | | | | | | Indene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Atrazine | 50 ug/mL |
| | | | | | | | Benzaldehyde | 50 ug/mL |
| | | | | | | | Caprolactam | 50 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 50 ug/mL |
| | | | | | | | Benzidine | 50 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 25 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 25 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 25 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 25 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 25 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 25 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 SS W_00005** | 04/20/16 | 04/21/15 | MECL2, Lot 0000093871 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| **SMLIST1 SS W_00005** | 04/20/16 | 04/21/15 | MECL2, Lot 0000093871 | 2 mL | SMLIST1 SS ST_00005 | 200 uL | 1,1'-Biphenyl | 10 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 10 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 10 ug/mL |
| | | | | | | | 2-Chlorophenol | 10 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 10 ug/mL |
| | | | | | | | 2-Methylphenol | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Nitroaniline | 10 ug/mL |
| | | | | | | | 2-Nitrophenol | 10 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 10 ug/mL |
| | | | | | | | 3-Nitroaniline | 10 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 20 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 10 ug/mL |
| | | | | | | | 4-Chloroaniline | 10 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Nitroaniline | 10 ug/mL |
| | | | | | | | 4-Nitrophenol | 20 ug/mL |
| | | | | | | | Acenaphthene | 10 ug/mL |
| | | | | | | | Acenaphthylene | 10 ug/mL |
| | | | | | | | Acetophenone | 10 ug/mL |
| | | | | | | | Anthracene | 10 ug/mL |
| | | | | | | | Benzo[a]anthracene | 10 ug/mL |
| | | | | | | | Benzo[a]pyrene | 10 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 10 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 10 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 10 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 10 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 10 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 10 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 10 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 10 ug/mL |
| | | | | | | | Carbazole | 10 ug/mL |
| | | | | | | | Chrysene | 10 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 10 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 10 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 10 ug/mL |
| | | | | | | | Dibenzofuran | 10 ug/mL |
| | | | | | | | Diethyl phthalate | 10 ug/mL |
| | | | | | | | Dimethyl phthalate | 10 ug/mL |
| | | | | | | | Fluoranthene | 10 ug/mL |
| | | | | | | | Fluorene | 10 ug/mL |
| | | | | | | | Hexachlorobenzene | 10 ug/mL |
| | | | | | | | Hexachlorobutadiene | 10 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 10 ug/mL |
| | | | | | | | Hexachloroethane | 10 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 10 ug/mL |
| | | | | | | | Isophorone | 10 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 10 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 20 ug/mL |
| | | | | | | | Naphthalene | 10 ug/mL |
| | | | | | | | Nitrobenzene | 10 ug/mL |
| | | | | | | | Pentachlorophenol | 20 ug/mL |
| | | | | | | | Phenanthrene | 10 ug/mL |
| | | | | | | | Phenol | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Pyrene | 10 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 20 ug/mL |
| | | | | | | | Atrazine | 20 ug/mL |
| | | | | | | | Benzaldehyde | 20 ug/mL |
| | | | | | | | Caprolactam | 20 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 10 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 10 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 10 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 10 ug/mL |
| .SMLIST1 SS ST_00005 | 06/30/16 | 04/21/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 SS S1_00004 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | SMLIST1 SS S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | SMLIST1SS S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1SS SUR_00002 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 SS S1_00004 | 06/30/16 | | Restek, Lot A0107697 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| ..SMLIST1 SS S9_00001 | 07/31/16 | | Restek, Lot A0108170 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..SMLIST1SS S11_00001 | 06/30/16 | | Restek, Lot A0108042 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1SS SUR_00002 | 09/30/18 | | Restek, Lot A092712 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **VGGASCALNEW_00014** | 09/22/15 | 06/23/15 | MEOH, Lot 62345 | 25 mL | VGGASSTOCKNEW_00003 | 1 mL | C6-C12 | 100 ug/mL |
| | | | | | | | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGGASSTOCKNEW_00003 | 01/31/20 | | RESTEK, Lot A092645 | | (Purchased Reagent) | | C6-C12 | 2500 ug/mL |
| | | | | | | | Gasoline Range Organics [C6 - C10] | 2500 ug/mL |
| **VGGASCALNEW_00015** | 12/23/15 | 09/23/15 | MEOH, Lot 62345 | 25 mL | VGGASSTOCKNEW_00003 | 1 mL | C6-C12 | 100 ug/mL |
| | | | | | | | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGGASSTOCKNEW_00003 | 01/31/20 | | RESTEK, Lot A092645 | | (Purchased Reagent) | | C6-C12 | 2500 ug/mL |
| | | | | | | | Gasoline Range Organics [C6 - C10] | 2500 ug/mL |
| **VGGASCHK_00068** | 10/27/15 | 07/27/15 | MEOH, Lot 49909 | 50 mL | VGASCHK_00002 | 1 mL | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGASCHK_00002 | 10/30/19 | | AccuStandard, Lot B5040009-01 | | (Purchased Reagent) | | Gasoline Range Organics [C6 - C10] | 5 mg/mL |
| **VGGASCHK_00070** | 02/04/16 | 11/04/15 | MEOH, Lot 49909 | 50 mL | VGASCHK_00002 | 1 mL | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGASCHK_00002 | 10/30/19 | | AccuStandard, Lot B5040009-01 | | (Purchased Reagent) | | Gasoline Range Organics [C6 - C10] | 5 mg/mL |
| **VGGASURROGATE_00034** | 01/09/16 | 10/09/15 | MEOH, Lot 49909 | 50 mL | VGSURRSTKGAS_00019 | 1.25 mL | Trifluorotoluene (Surr) | 50 ug/mL |
| .VGSURRSTKGAS_00019 | 08/31/16 | | ULTRA SCIENTIFIC, Lot CE-1903A | | (Purchased Reagent) | | Trifluorotoluene (Surr) | 2000 ug/mL |
| **VM250-IS_00028** | 09/13/15 | 03/13/15 | MEOH, Lot 0000085233 | 100 mL | vmcus8126_00008 | 1 mL | 1,4-Dichlorobenzene-d4 | 250 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 250 ug/mL |
| | | | | | | | Fluorobenzene | 250 ug/mL |
| .vmcus8126_00008 | 09/30/16 | | Ultra Scientific, Lot CL-3192 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 25000 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 25000 ug/mL |
| | | | | | | | Fluorobenzene | 25000 ug/mL |
| **VM250-IS_00029** | 10/10/15 | 04/10/15 | MEOH, Lot 0000085233 | 100 mL | vmcus8126_00008 | 1 mL | 1,4-Dichlorobenzene-d4 | 250 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 250 ug/mL |
| | | | | | | | Fluorobenzene | 250 ug/mL |
| .vmcus8126_00008 | 09/30/16 | | Ultra Scientific, Lot CL-3192 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 25000 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 25000 ug/mL |
| | | | | | | | Fluorobenzene | 25000 ug/mL |
| **VM250-IS_00030** | 03/17/16 | 09/17/15 | MEOH, Lot +1632000000994941 | 100 mL | vmcus8126_00008 | 1 mL | 1,4-Dichlorobenzene-d4 | 250 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 250 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Fluorobenzene | 250 ug/mL |
| .vmcus8126_00008 | 09/30/16 | Ultra Scientific, Lot CL-3192 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 25000 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 25000 ug/mL |
| | | | | | | | Fluorobenzene | 25000 ug/mL |
| **VM250SS_00031** | 11/29/15 | 05/29/15 | MEOH, Lot +161620000085233JB | 50 mL | VM567650_00021 | 5 mL | 1,2-Dichloroethane-d4 (Surr) | 250 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 250 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 250 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 250 ug/mL |
| .VM567650_00021 | 04/30/19 | Restek, Lot A0102817 | | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **VM250SS_00033** | 02/09/16 | 11/09/15 | MEOH, Lot +171940000118655JM | 50 mL | VM567650_00023 | 5 mL | 1,2-Dichloroethane-d4 (Surr) | 250 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 250 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 250 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 250 ug/mL |
| .VM567650_00023 | 06/30/19 | Restek, Lot A0104073 | | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **VM50IS_00050** | 12/11/15 | 06/11/15 | MEOH, Lot 85233 | 100 mL | vm30241_00002 | 2 mL | 1,4-Dichlorobenzene-d4 | 50 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 50 ug/mL |
| | | | | | | | Fluorobenzene | 50 ug/mL |
| .vm30241_00002 | 11/30/19 | restek, Lot A0107133 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2500 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 2500 ug/mL |
| | | | | | | | Fluorobenzene | 2500 ug/mL |
| **VM50IS_00052** | 04/15/16 | 10/15/15 | MEOH, Lot 99494 | 100 mL | vm30241_00002 | 2 mL | 1,4-Dichlorobenzene-d4 | 50 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 50 ug/mL |
| | | | | | | | Fluorobenzene | 50 ug/mL |
| .vm30241_00002 | 11/30/19 | restek, Lot A0107133 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2500 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 2500 ug/mL |
| | | | | | | | Fluorobenzene | 2500 ug/mL |
| **vm50ss_00210** | 08/05/15 | 07/29/15 | MEOH, Lot na | 2 mL | vm50ss_stk_00065 | 2 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| .vm50ss_stk_00065 | 09/08/15 | 03/08/15 | MEOH, Lot 0000085233 | 200 mL | VM567650_00021 | 4 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| ..VM567650_00021 | 04/30/19 | Restek, Lot A0102817 | | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **vm50ss_00213** | 09/02/15 | 08/26/15 | MEOH, Lot na | 2 mL | vm50ss_stk_00065 | 2 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| .vm50ss_stk_00065 | 09/08/15 | 03/08/15 | MEOH, Lot 0000085233 | 200 mL | VM567650_00021 | 4 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| ..VM567650_00021 | 04/30/19 | | Restek, Lot A0102817 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **vm50ss_00220** | 11/23/15 | 11/16/15 | MEOH, Lot na | 2 mL | vm50ss_stk_00067 | 2 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| .vm50ss_stk_00067 | 04/05/16 | 10/05/15 | MEOH, Lot 0000099494 | 200 mL | VM567650_00023 | 4 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| ..VM567650_00023 | 06/30/19 | | Restek, Lot A0104073 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **vm50ss_stk_00067** | 04/05/16 | 10/05/15 | MEOH, Lot 0000099494 | 200 mL | VM567650_00023 | 4 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| .VM567650_00023 | 06/30/19 | | Restek, Lot A0104073 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **VMAROLISTDW_00109** | 07/31/15 | 07/24/15 | MEOH, Lot na | 3 mL | VMACROLSTD_00027 | 3 mL | Acrolein | 250 ug/mL |
| .VMACROLSTD_00027 | 07/31/15 | 07/08/15 | MEOH, Lot +161620000085233JB | 20 mL | VM568720_00007 | 0.25 mL | Acrolein | 250 ug/mL |
| ..VM568720_00007 | 07/31/15 | | restek, Lot A0109948 | | (Purchased Reagent) | | Acrolein | 20000 ug/mL |
| **VMAROLISTDW_00114** | 09/08/15 | 09/01/15 | MEOH, Lot na | 3 mL | VMACROLSTD_00029 | 3 mL | Acrolein | 250 ug/mL |
| .VMACROLSTD_00029 | 09/30/15 | 09/01/15 | MEOH, Lot 147358 | 20 mL | VM568720_00008 | 250 uL | Acrolein | 250 ug/mL |
| ..VM568720_00008 | 09/30/15 | | restek, Lot A0111006 | | (Purchased Reagent) | | Acrolein | 20000 ug/mL |
| **VMAROLISTDW_00125** | 11/23/15 | 11/16/15 | MEOH, Lot na | 3 mL | VMACROLSTD_00031 | 3 mL | Acrolein | 250 ug/mL |
| .VMACROLSTD_00031 | 11/30/15 | 11/01/15 | MEOH, Lot 0000118655 | 20 mL | VM568720_00009 | 250 uL | Acrolein | 250 ug/mL |
| ..VM568720_00009 | 11/30/15 | | restek, Lot A0112264 | | (Purchased Reagent) | | Acrolein | 20000 ug/mL |
| **VMFASA9W_00077** | 11/23/15 | 11/16/15 | MEOH, Lot NA | 2 mL | VMFASA9_00005 | 2 mL | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| .VMFASA9_00005 | 11/30/15 | 06/30/15 | MEOH, Lot +161620000085233jb | 100 mL | VM568723S_00001 | 2.5 mL | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725S_00001 | 2 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| ..VM568723S_00001 | 04/30/16 | restek, Lot A0102749 | | | (Purchased Reagent) | | Diisopropyl ether | 2000 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725S_00001 | 11/30/15 | restek, Lot A0111261 | | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| **VMFASG_00033** | 09/29/15 | 08/29/15 | MEOH, Lot 147358 | 50 mL | VM567645S_00016 | 1.25 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .VM567645S_00016 | 11/30/15 | Restek, Lot A099261 | | | (Purchased Reagent) | | Bromomethane | 2000 ug/mL |
| | | | | | | | Chloroethane | 2000 ug/mL |
| | | | | | | | Chloromethane | 2000 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2000 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2000 ug/mL |
| | | | | | | | Vinyl chloride | 2000 ug/mL |
| **VMFASGW_00123** | 11/19/15 | 11/12/15 | MEOH, Lot NA | 2 mL | VMFASG_00035 | 2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .VMFASG_00035 | 11/30/15 | 10/27/15 | MEOH, Lot 0000118655 | 50 mL | VM567645S_00016 | 1.25 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| ..VM567645S_00016 | 11/30/15 | Restek, Lot A099261 | | | (Purchased Reagent) | | Bromomethane | 2000 ug/mL |
| | | | | | | | Chloroethane | 2000 ug/mL |
| | | | | | | | Chloromethane | 2000 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2000 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2000 ug/mL |
| | | | | | | | Vinyl chloride | 2000 ug/mL |
| **VMFASGW_00124** | 11/23/15 | 11/16/15 | MEOH, Lot NA | 2 mL | VMFASG_00036 | 2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .VMFASG_00036 | 11/30/15 | 11/16/15 | MEOH, Lot 0000118655 | 50 mL | VM567645S_00016 | 1.25 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| ..VM567645S_00016 | 11/30/15 | | Restek, Lot A099261 | | (Purchased Reagent) | | Bromomethane | 2000 ug/mL |
| | | | | | | | Chloroethane | 2000 ug/mL |
| | | | | | | | Chloromethane | 2000 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2000 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2000 ug/mL |
| | | | | | | | Vinyl chloride | 2000 ug/mL |
| **VMFASPW_00114** | 09/08/15 | 09/01/15 | MEOH, Lot n/a | 2 mL | VMRFASP_00022 | 2 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRFASP_00022 | 10/31/15 | 07/28/15 | MEOH, Lot 147358 | 100 mL | VM567641S_00007 | 2.5 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | VM567642S_00005 | 1 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM569723S_00002 | 2 mL | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | VM569724S_00002 | 1 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641S_00007 | 02/29/16 | | Restek, Lot A093733 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Xylenes, Total | 4000 ug/mL |
| ..VM567642S_00005 | 02/29/16 | | Restek, Lot A093472 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM569723S_00002 | 03/31/18 | | Restek, Lot A0110013 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2500 ug/mL |
| ..VM569724S_00002 | 10/31/15 | | Restek, Lot A0110474 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMFASPW_00126** | 11/19/15 | 11/12/15 | MEOH, Lot n/a | 2 mL | VMRFASP_00024 | 2 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRFASP_00024 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 100 mL | VM567641S_00007 | 2.5 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | VM567642S_00005 | 1 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM569723S_00001 | 2 mL | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | VM569724S_00003 | 1 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641S_00007 | 02/29/16 | | Restek, Lot A093733 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| | | | | | | | Xylenes, Total | 4000 ug/mL |
| ..VM567642S_00005 | 02/29/16 | | Restek, Lot A093472 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM569723S_00001 | 01/31/18 | | Restek, Lot A0108188 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2500 ug/mL |
| ..VM569724S_00003 | 01/31/16 | | Restek, Lot A0112364 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMFASPW_00127** | 11/27/15 | 11/20/15 | MEOH, Lot n/a | 2 mL | VMRFASP_00024 | 2 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRFASP_00024 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 100 mL | VM567641S_00007 | 2.5 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | VM567642S_00005 | 1 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM569723S_00001 | 2 mL | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | VM569724S_00003 | 1 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641S_00007 | 02/29/16 | | Restek, Lot A093733 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| | | | | | | | Xylenes, Total | 4000 ug/mL |
| ..VM567642S_00005 | 02/29/16 | | Restek, Lot A093472 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM569723S_00001 | 01/31/18 | | Restek, Lot A0108188 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..VM569724S_00003 | 01/31/16 | Restek, Lot A0112364 | | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRA9W_00115** | 07/22/15 | 07/15/15 | MEOH, Lot NA | 1 mL | VMRA9_00013 | 1 mL | Cyclohexanone | 500 ug/mL |
| | | | | | | | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| .VMRA9_00013 | 11/30/15 | 06/30/15 | MEOH, Lot +161620000085233jb | 50 mL | VM567648_00014 | 1.25 mL | Cyclohexanone | 500 ug/mL |
| | | | | | VM568723_00001 | 1.25 mL | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725_00002 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | vm569726_00001 | 2 mL | 2-Methylnaphthalene | 100 ug/mL |
| ..VM567648_00014 | 02/29/16 | Restek, Lot A093361 | | | (Purchased Reagent) | | Cyclohexanone | 20000 ug/mL |
| ..VM568723_00001 | 12/31/15 | restek, Lot A099930 | | | (Purchased Reagent) | | Acetonitrile | 20000 ug/mL |
| | | | | | | | Diisopropyl ether | 2000 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Propionitrile | 20000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725_00002 | 11/30/15 | restek, Lot A0111228 | | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 2500 ug/mL |
| | | | | | | | 1-Chlorohexane | 2500 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 2500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Nitropropane | 5000 ug/mL |
| | | | | | | | Ethyl acetate | 5000 ug/mL |
| | | | | | | | Ethyl acrylate | 2500 ug/mL |
| | | | | | | | Methacrylonitrile | 25000 ug/mL |
| | | | | | | | Methyl methacrylate | 5000 ug/mL |
| | | | | | | | n-Butanol | 62500 ug/mL |
| | | | | | | | n-Butyl acetate | 2500 ug/mL |
| ..vm569726_00001 | 04/30/17 | | Restek, Lot A0110130 | | (Purchased Reagent) | | 2-Methylnaphthalene | 2500 ug/mL |
| **VMRA9W_00119** | 08/17/15 | 08/10/15 | MEOH, Lot NA | 1 mL | VMRA9_00013 | 1 mL | Cyclohexanone | 500 ug/mL |
| | | | | | | | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| .VMRA9_00013 | 11/30/15 | 06/30/15 | MEOH, Lot +161620000085233jb | 50 mL | VM567648_00014 | 1.25 mL | Cyclohexanone | 500 ug/mL |
| | | | | | VM568723_00001 | 1.25 mL | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725_00002 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | vm569726_00001 | 2 mL | 2-Methylnaphthalene | 100 ug/mL |
| ..VM567648_00014 | 02/29/16 | | Restek, Lot A093361 | | (Purchased Reagent) | | Cyclohexanone | 20000 ug/mL |
| ..VM568723_00001 | 12/31/15 | | restek, Lot A099930 | | (Purchased Reagent) | | Acetonitrile | 20000 ug/mL |
| | | | | | | | Diisopropyl ether | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Propionitrile | 20000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725_00002 | 11/30/15 | | restek, Lot A0111228 | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 2500 ug/mL |
| | | | | | | | 1-Chlorohexane | 2500 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 2500 ug/mL |
| | | | | | | | 2-Nitropropane | 5000 ug/mL |
| | | | | | | | Ethyl acetate | 5000 ug/mL |
| | | | | | | | Ethyl acrylate | 2500 ug/mL |
| | | | | | | | Methacrylonitrile | 25000 ug/mL |
| | | | | | | | Methyl methacrylate | 5000 ug/mL |
| | | | | | | | n-Butanol | 62500 ug/mL |
| | | | | | | | n-Butyl acetate | 2500 ug/mL |
| ..vm569726_00001 | 04/30/17 | | Restek, Lot A0110130 | | (Purchased Reagent) | | 2-Methylnaphthalene | 2500 ug/mL |
| **VMRA9W_00131** | 11/23/15 | 11/16/15 | MEOH, Lot NA | 1 mL | VMRA9_00013 | 1 mL | Cyclohexanone | 500 ug/mL |
| | | | | | | | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| .VMRA9_00013 | 11/30/15 | 06/30/15 | MEOH, Lot +161620000085233jb | 50 mL | VM567648_00014 | 1.25 mL | Cyclohexanone | 500 ug/mL |
| | | | | | VM568723_00001 | 1.25 mL | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725_00002 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | vm569726_00001 | 2 mL | 2-Methylnaphthalene | 100 ug/mL |
| ..VM567648_00014 | 02/29/16 | | Restek, Lot A093361 | | (Purchased Reagent) | | Cyclohexanone | 20000 ug/mL |
| ..VM568723_00001 | 12/31/15 | | restek, Lot A099930 | | (Purchased Reagent) | | Acetonitrile | 20000 ug/mL |
| | | | | | | | Diisopropyl ether | 2000 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Propionitrile | 20000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725_00002 | 11/30/15 | | restek, Lot A0111228 | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 2500 ug/mL |
| | | | | | | | 1-Chlorohexane | 2500 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 2500 ug/mL |
| | | | | | | | 2-Nitropropane | 5000 ug/mL |
| | | | | | | | Ethyl acetate | 5000 ug/mL |
| | | | | | | | Ethyl acrylate | 2500 ug/mL |
| | | | | | | | Methacrylonitrile | 25000 ug/mL |
| | | | | | | | Methyl methacrylate | 5000 ug/mL |
| | | | | | | | n-Butanol | 62500 ug/mL |
| | | | | | | | n-Butyl acetate | 2500 ug/mL |
| ..vm569726_00001 | 04/30/17 | | Restek, Lot A0110130 | | (Purchased Reagent) | | 2-Methylnaphthalene | 2500 ug/mL |
| **VMRGAS_00112** | 08/04/15 | 07/28/15 | MEOH, Lot 147358 | 10 mL | vm569722_00001 | 0.2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Butadiene | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Dichlorofluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .vm569722_00001 | 01/31/18 | | Restek, Lot A0108198 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Butadiene | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Dichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMRGAS_00116** | 09/05/15 | 08/29/15 | MEOH, Lot 147358 | 10 mL | vm569722_00002 | 0.2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Butadiene | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Dichlorofluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .vm569722_00002 | 01/31/18 | | Restek, Lot A0108198 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Butadiene | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Dichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMRGAS_00126** | 11/20/15 | 11/13/15 | MEOH, Lot 0000118655 | 10 mL | vm569722_00003 | 0.2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Butadiene | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Dichlorofluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .vm569722_00003 | 04/30/18 | | Restek, Lot A0110070 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Butadiene | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Dichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMRPRIMW_00140** | 07/31/15 | 07/24/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00012 | 1 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRPRIM_00012 | 08/31/15 | 05/27/15 | MEOH, Lot +161620000085233JB | 50 mL | VM567641_00013 | 1.25 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | VM567642_00015 | 0.5 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM567643_00020 | 2.5 mL | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | VM569724_00001 | 0.5 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641_00013 | 02/29/16 | | restek, Lot A093581 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,4-Dioxane | 40000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Acrylonitrile | 20000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Cyclohexane | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethyl ether | 2000 ug/mL |
| | | | | | | | Ethyl methacrylate | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Hexane | 2000 ug/mL |
| | | | | | | | Iodomethane | 2000 ug/mL |
| | | | | | | | Isobutyl alcohol | 50000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl acetate | 10000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylcyclohexane | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | n-Heptane | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Tetrahydrofuran | 4000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| ..VM567642_00015 | 02/29/16 | | Restek, Lot A093365 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM567643_00020 | 02/29/16 | | restek, Lot A093368 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2000 ug/mL |
| ..VM569724_00001 | 08/31/15 | | Restek, Lot A0109190 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRPRIMW_00146** | 09/08/15 | 09/01/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00013 | 1 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRPRIM_00013 | 10/31/15 | 09/01/15 | MEOH, Lot 147358 | 50 mL | VM567641_00013 | 1.25 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | VM567642_00015 | 0.5 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM567643_00020 | 2.5 mL | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | VM569724_00002 | 0.5 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641_00013 | 02/29/16 | | restek, Lot A093581 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,4-Dioxane | 40000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Acrylonitrile | 20000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Cyclohexane | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethyl ether | 2000 ug/mL |
| | | | | | | | Ethyl methacrylate | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Hexane | 2000 ug/mL |
| | | | | | | | Iodomethane | 2000 ug/mL |
| | | | | | | | Isobutyl alcohol | 50000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl acetate | 10000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylcyclohexane | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | n-Heptane | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Tetrahydrofuran | 4000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| ..VM567642_00015 | 02/29/16 | | Restek, Lot A093365 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM567643_00020 | 02/29/16 | | restek, Lot A093368 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2000 ug/mL |
| ..VM569724_00002 | 10/31/15 | | Restek, Lot A0110407 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRPRIMW_00157** | 11/23/15 | 11/16/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00014 | 1 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRPRIM_00014 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 50 mL | VM567641_00013 | 1.25 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | VM567642_00015 | 0.5 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM567643_00020 | 2.5 mL | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | VM569724_00003 | 0.5 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641_00013 | 02/29/16 | | restek, Lot A093581 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,4-Dioxane | 40000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Acrylonitrile | 20000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Cyclohexane | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethyl ether | 2000 ug/mL |
| | | | | | | | Ethyl methacrylate | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Hexane | 2000 ug/mL |
| | | | | | | | Iodomethane | 2000 ug/mL |
| | | | | | | | Isobutyl alcohol | 50000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl acetate | 10000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylcyclohexane | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | n-Heptane | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Tetrahydrofuran | 4000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| ..VM567642_00015 | 02/29/16 | Restek, Lot A093365 | | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM567643_00020 | 02/29/16 | restek, Lot A093368 | | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2000 ug/mL |
| ..VM569724_00003 | 01/31/16 | Restek, Lot A0112975 | | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRPRIMW_00157** | 11/23/15 | 11/16/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00014 | 1 mL | Xylenes, Total | 100 ug/mL |
| .VMRPRIM_00014 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 50 mL | VM567641_00013 | 1.25 mL | Xylenes, Total | 100 ug/mL |
| ..VM567641_00013 | 02/29/16 | restek, Lot A093581 | | | (Purchased Reagent) | | Xylenes, Total | 4000 ug/mL |
| **VMUX8SS_00022** | 01/11/16 | 10/11/15 | MEOH, Lot 0000099494 | 100 mL | VMCUS8125_00009 | 866 uL | 1,2-Dichloroethane-d4 (Surr) | 216.5 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 216.5 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 216.5 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 216.5 ug/mL |
| .VMCUS8125_00009 | 09/30/16 | Ultra Scientific, Lot CJ-3208 | | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 25000 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 25000 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 25000 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 25000 ug/mL |

# Certification Summary

Client: Tetra Tech, Inc. TestAmerica Job ID: 240-57930-1
Project/Site: LM MRC Block F UST Pipe Inv.

| Laboratory | Authority | Program | EPA Region | Certification ID |
|---|---|---|---|---|
| TestAmerica Canton | California | State Program | 9 | 2927 |
| TestAmerica Canton | Connecticut | State Program | 1 | PH-0590 |
| TestAmerica Canton | Illinois | NELAP | 5 | 200004 |
| TestAmerica Canton | Kansas | NELAP | 7 | E-10336 |
| TestAmerica Canton | Kentucky (UST) | State Program | 4 | 58 |
| TestAmerica Canton | Kentucky (WW) | State Program | 4 | 98016 |
| TestAmerica Canton | L-A-B | DoD ELAP | | L2315 |
| TestAmerica Canton | Minnesota | NELAP | 5 | 039-999-348 |
| TestAmerica Canton | Nevada | State Program | 9 | OH-000482008A |
| TestAmerica Canton | New Jersey | NELAP | 2 | OH001 |
| TestAmerica Canton | New York | NELAP | 2 | 10975 |
| TestAmerica Canton | Ohio VAP | State Program | 5 | CL0024 |
| TestAmerica Canton | Oregon | NELAP | 10 | 4062 |
| TestAmerica Canton | Pennsylvania | NELAP | 3 | 68-00340 |
| TestAmerica Canton | Texas | NELAP | 6 | T104704517-15-5 |
| TestAmerica Canton | USDA | Federal | | P330-13-00319 |
| TestAmerica Canton | Virginia | NELAP | 3 | 460175 |
| TestAmerica Canton | Washington | State Program | 10 | C971 |
| TestAmerica Canton | West Virginia DEP | State Program | 3 | 210 |
| TestAmerica Canton | Wisconsin | State Program | 5 | 999518190 |

Accreditation may not be offered or required for all methods and analytes reported in this package. Please contact your project manager for the laboratory's current list of certified methods and analytes.

# Method 8260C

## Volatile Organic Compounds (GC/MS) by Method 8260C

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): DB-624 ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | DBFM # | DCA # | TOL # | BFB # |
|---|---|---|---|---|---|
| F-VS-UST-P-1 | 240-57930-1 | 100 | 104 | 104 | 97 |
| F-VS-UST-P-2 | 240-57930-2 | 92 | 98 | 94 | 91 |
| F-VS-UST-P-2 | 240-57930-2 | 92 | 95 | 103 | 105 |
| F-VS-UST-P-2 | 240-57930-2 | 94 | 88 | 110 | 93 |
| F-VS-UST-P-3 | 240-57930-3 | 91 | 91 | 94 | 89 |
| F-VS-UST-P-4 | 240-57930-4 | 87 | 85 | 111 | 102 * |
| F-VB-UST-P-1 | 240-57930-5 | 89 | 90 | 89 | 78 |
| F-VB-UST-P-2 | 240-57930-6 | 95 | 100 | 89 | 79 |
| F-VB-UST-P-3 | 240-57930-7 | 99 | 105 | 94 | 82 |
| F-VB-UST-P-4 | 240-57930-8 | 103 | 113 | 96 | 89 |
| F-VB-UST-P-5 | 240-57930-9 | 86 | 89 | 84 | 76 |
| F-VB-UST-P-6 | 240-57930-10 | 97 | 105 | 93 | 86 |
| F-VS-UST-P-DUP | 240-57930-11 | 89 | 94 | 90 | 83 |
| | MB 240-207343/6 | 91 | 95 | 87 | 76 |
| | MB 240-207560/6 | 98 | 102 | 92 | 82 |
| | MB 240-207720/8 | 117 | 114 | 120 | 114 |
| | LCS 240-207343/5 | 97 | 94 | 91 | 80 |
| | LCS 240-207343/7 | 90 | 92 | 89 | 80 |
| | LCS 240-207560/5 | 98 | 98 | 92 | 81 |
| | LCS 240-207560/7 | 95 | 94 | 89 | 83 |
| | LCS 240-207720/6 | 105 | 100 | 104 | 96 |
| | 240-57921-A-4-J MS | 52 | 93 | 90 | 83 |
| | 240-57972-C-3-C MS | 101 | 94 | 101 | 98 |
| | 240-57921-A-4-K MSD | 49 | 93 | 90 | 84 |
| | 240-57972-C-3-D MSD | 98 | 95 | 101 | 100 |

| | QC LIMITS |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 37-132 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 58-123 |
| TOL = Toluene-d8 (Surr) | 67-125 |
| BFB = 4-Bromofluorobenzene (Surr) | 52-136 |

# Column to be used to flag recovery values

FORM II 8260C

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Water Level: Low

GC Column (1): DB-624 ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | DBFM # | DCA # | TOL # | BFB # |
|---|---|---|---|---|---|
| TB-111315 | 240-57930-12 | 81 | 83 | 82 | 75 |
| | MB 240-207096/6 | 82 | 82 | 82 | 75 |
| | LCS 240-207096/4 | 90 | 90 | 90 | 85 |
| | 200-30727-B-1 MS | 81 | 81 | 84 | 80 |
| | 200-30727-B-1 MSD | 84 | 80 | 84 | 80 |

| | QC LIMITS |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 79-120 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 78-125 |
| TOL = Toluene-d8 (Surr) | 80-120 |
| BFB = 4-Bromofluorobenzene (Surr) | 61-120 |

# Column to be used to flag recovery values

FORM II 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXX3068.D

Lab ID: LCS 240-207096/4 Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 10.0 | 9.46 | 95 | 80-120 | |
| 1,1,1-Trichloroethane | 10.0 | 9.11 | 91 | 77-123 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 10.0 | 100 | 71-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 10.0 | 9.72 | 97 | 67-138 | |
| 1,1-Dichloroethane | 10.0 | 9.27 | 93 | 79-125 | |
| 1,1-Dichloroethene | 10.0 | 9.08 | 91 | 76-124 | |
| 1,1-Dichloropropene | 10.0 | 10.0 | 100 | 80-128 | |
| 1,2,3-Trichlorobenzene | 10.0 | 5.89 | 59 | 66-120 | * |
| 1,2,3-Trichloropropane | 10.0 | 10.0 | 100 | 72-126 | |
| 1,2,4-Trichlorobenzene | 10.0 | 6.37 | 64 | 61-120 | |
| 1,2,4-Trimethylbenzene | 10.0 | 9.26 | 93 | 76-120 | |
| 1,2-Dibromo-3-Chloropropane | 10.0 | 8.37 | 84 | 50-132 | |
| 1,2-Dichlorobenzene | 10.0 | 9.01 | 90 | 79-120 | |
| 1,2-Dichloroethane | 10.0 | 9.37 | 94 | 80-120 | |
| 1,2-Dichloropropane | 10.0 | 9.44 | 94 | 78-124 | |
| 1,3-Dichlorobenzene | 10.0 | 9.15 | 91 | 79-120 | |
| 1,3-Dichloropropane | 10.0 | 9.46 | 95 | 80-120 | |
| 1,4-Dichlorobenzene | 10.0 | 9.15 | 91 | 79-120 | |
| 2,2-Dichloropropane | 10.0 | 7.77 | 78 | 71-122 | |
| 2-Butanone (MEK) | 20.0 | 19.0 | 95 | 56-138 | |
| 2-Chloroethyl vinyl ether | 10.0 | 9.41 J | 94 | 54-139 | |
| 2-Chlorotoluene | 10.0 | 9.78 | 98 | 80-120 | |
| 2-Hexanone | 20.0 | 19.2 | 96 | 55-141 | |
| 4-Chlorotoluene | 10.0 | 9.88 | 99 | 80-120 | |
| p-Isopropyltoluene | 10.0 | 9.20 | 92 | 77-120 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 18.8 | 94 | 64-135 | |
| Acetone | 20.0 | 18.4 | 92 | 34-148 | |
| Benzene | 10.0 | 9.16 | 92 | 80-120 | |
| Bromobenzene | 10.0 | 9.75 | 98 | 80-120 | |
| Bromoform | 10.0 | 8.92 | 89 | 56-122 | |
| Bromomethane | 10.0 | 8.49 | 85 | 38-132 | |
| Carbon disulfide | 10.0 | 8.54 | 85 | 65-144 | |
| Carbon tetrachloride | 10.0 | 9.28 | 93 | 77-131 | |
| Chlorobenzene | 10.0 | 9.38 | 94 | 80-120 | |
| Dibromochloromethane | 10.0 | 9.64 | 96 | 74-120 | |
| Chloroethane | 10.0 | 8.47 | 85 | 36-126 | |
| Chloroform | 10.0 | 9.47 | 95 | 80-120 | |
| Chloromethane | 10.0 | 8.43 | 84 | 48-133 | |
| cis-1,2-Dichloroethene | 10.0 | 9.25 | 92 | 79-120 | |
| cis-1,3-Dichloropropene | 10.0 | 9.55 | 96 | 74-126 | |
| Dibromomethane | 10.0 | 9.34 | 93 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXX3068.D

Lab ID: LCS 240-207096/4 Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 10.0 | 9.25 | 93 | 80-120 | |
| Dichlorodifluoromethane | 10.0 | 6.79 | 68 | 23-136 | |
| Bromodichloromethane | 10.0 | 9.05 | 90 | 80-120 | |
| tert-Butyl alcohol (TBA) | 100 | 73.6 | 74 | 38-144 | |
| Ethylbenzene | 10.0 | 9.52 | 95 | 80-120 | |
| Hexachlorobutadiene | 10.0 | 6.28 | 63 | 62-120 | |
| Isopropylbenzene | 10.0 | 9.49 | 95 | 77-120 | |
| Methyl tert-butyl ether | 10.0 | 8.83 | 88 | 69-121 | |
| Methylene Chloride | 10.0 | 10.9 | 109 | 77-129 | |
| m-Xylene & p-Xylene | 10.0 | 9.72 | 97 | 80-120 | |
| Naphthalene | 10.0 | 6.29 | 63 | 53-121 | |
| n-Butylbenzene | 10.0 | 8.80 | 88 | 74-120 | |
| N-Propylbenzene | 10.0 | 10.1 | 101 | 79-123 | |
| o-Xylene | 10.0 | 9.71 | 97 | 80-120 | |
| sec-Butylbenzene | 10.0 | 9.36 | 94 | 77-120 | |
| Styrene | 10.0 | 9.41 | 94 | 76-122 | |
| Tetrachloroethene | 10.0 | 9.47 | 95 | 78-121 | |
| Toluene | 10.0 | 9.21 | 92 | 80-120 | |
| trans-1,2-Dichloroethene | 10.0 | 9.18 | 92 | 80-124 | |
| trans-1,3-Dichloropropene | 10.0 | 9.95 | 100 | 75-131 | |
| Trichloroethene | 10.0 | 9.58 | 96 | 80-121 | |
| Trichlorofluoromethane | 10.0 | 9.63 | 96 | 61-133 | |
| Vinyl acetate | 10.0 | 7.50 | 75 | 27-160 | |
| Vinyl chloride | 10.0 | 8.72 | 87 | 52-121 | |
| Xylenes, Total | 20.0 | 19.4 | 97 | 80-120 | |
| tert-Butylbenzene | 10.0 | 9.70 | 97 | 77-121 | |
| 1,2-Dibromoethane | 10.0 | 9.91 | 99 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146534.D

Lab ID: LCS 240-207343/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 21.0 | 105 | 80-120 | |
| 1,1,1-Trichloroethane | 20.0 | 22.4 | 112 | 77-126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 17.6 | 88 | 77-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 20.3 | 101 | 80-138 | |
| 1,1-Dichloroethane | 20.0 | 18.2 | 91 | 76-120 | |
| 1,1-Dichloroethene | 20.0 | 17.0 | 85 | 75-135 | |
| 1,1-Dichloropropene | 20.0 | 19.3 | 97 | 80-120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 18.7 | 93 | 61-121 | |
| 1,2,3-Trichloropropane | 20.0 | 20.0 | 100 | 80-126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 18.7 | 93 | 64-124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 18.7 | 94 | 80-129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 18.1 | 90 | 61-132 | |
| 1,2-Dichlorobenzene | 20.0 | 18.0 | 90 | 76-120 | |
| 1,2-Dichloroethane | 20.0 | 19.6 | 98 | 72-120 | |
| 1,2-Dichloropropane | 20.0 | 16.9 | 84 | 80-120 | |
| 1,3-Dichlorobenzene | 20.0 | 17.9 | 89 | 78-120 | |
| 1,3-Dichloropropane | 20.0 | 17.4 | 87 | 80-120 | |
| 1,4-Dichlorobenzene | 20.0 | 17.6 | 88 | 75-120 | |
| 2,2-Dichloropropane | 20.0 | 22.1 | 111 | 69-135 | |
| 2-Butanone (MEK) | 40.0 | 26.6 | 67 | 52-131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 12.5 J | 62 | 28-173 | |
| 2-Chlorotoluene | 20.0 | 19.3 | 97 | 78-120 | |
| 2-Hexanone | 40.0 | 30.6 | 76 | 64-136 | |
| 4-Chlorotoluene | 20.0 | 19.5 | 97 | 79-120 | |
| p-Isopropyltoluene | 20.0 | 19.3 | 97 | 77-131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 30.2 | 76 | 67-135 | |
| Acetone | 40.0 | 36.6 | 92 | 41-137 | |
| Benzene | 20.0 | 18.3 | 91 | 79-120 | |
| Bromobenzene | 20.0 | 19.7 | 99 | 80-120 | |
| Bromoform | 20.0 | 16.8 | 84 | 62-133 | |
| Bromomethane | 20.0 | 18.2 | 91 | 42-136 | |
| Carbon disulfide | 20.0 | 16.2 | 81 | 62-146 | |
| Carbon tetrachloride | 20.0 | 22.6 | 113 | 71-129 | |
| Chlorobenzene | 20.0 | 18.9 | 95 | 78-120 | |
| Dibromochloromethane | 20.0 | 18.3 | 91 | 72-127 | |
| Chloroethane | 20.0 | 17.0 | 85 | 58-120 | |
| Chloroform | 20.0 | 19.9 | 100 | 77-120 | |
| Chloromethane | 20.0 | 13.3 | 67 | 50-120 | |
| cis-1,2-Dichloroethene | 20.0 | 20.0 | 100 | 76-120 | |
| cis-1,3-Dichloropropene | 20.0 | 16.5 | 82 | 74-128 | |
| Dibromomethane | 20.0 | 19.1 | 95 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146534.D

Lab ID: LCS 240-207343/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 20.0 | 20.6 | 103 | 79-120 | |
| Dichlorodifluoromethane | 20.0 | 16.4 | 82 | 26-120 | |
| Bromodichloromethane | 20.0 | 18.4 | 92 | 80-122 | |
| tert-Butyl alcohol (TBA) | 200 | 135 J | 68 | 60-125 | |
| Ethylbenzene | 20.0 | 19.3 | 97 | 79-120 | |
| Hexachlorobutadiene | 20.0 | 19.5 | 97 | 54-131 | |
| Isopropylbenzene | 20.0 | 19.6 | 98 | 76-122 | |
| Methyl tert-butyl ether | 20.0 | 18.0 | 90 | 49-165 | |
| Methylene Chloride | 20.0 | 15.3 | 76 | 75-120 | |
| m-Xylene & p-Xylene | 20.0 | 18.8 | 94 | 80-120 | |
| Naphthalene | 20.0 | 16.8 | 84 | 65-123 | |
| n-Butylbenzene | 20.0 | 18.9 | 95 | 68-135 | |
| N-Propylbenzene | 20.0 | 19.0 | 95 | 80-129 | |
| o-Xylene | 20.0 | 20.1 | 101 | 80-120 | |
| sec-Butylbenzene | 20.0 | 18.6 | 93 | 74-129 | |
| Styrene | 20.0 | 18.2 | 91 | 80-120 | |
| Tetrachloroethene | 20.0 | 20.1 | 101 | 79-120 | |
| Toluene | 20.0 | 19.5 | 97 | 75-120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.1 | 100 | 78-120 | |
| trans-1,3-Dichloropropene | 20.0 | 18.1 | 90 | 73-131 | |
| Trichloroethene | 20.0 | 20.3 | 102 | 79-120 | |
| Trichlorofluoromethane | 20.0 | 20.8 | 104 | 57-146 | |
| Vinyl acetate | 20.0 | 15.5 | 77 | 70-130 | |
| Vinyl chloride | 20.0 | 14.2 | 71 | 57-120 | |
| Xylenes, Total | 40.0 | 38.9 | 97 | 80-120 | |
| tert-Butylbenzene | 20.0 | 19.0 | 95 | 76-126 | |
| 1,2-Dibromoethane | 20.0 | 20.0 | 100 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146567.D

Lab ID: LCS 240-207560/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 21.2 | 106 | 80-120 | |
| 1,1,1-Trichloroethane | 20.0 | 23.2 | 116 | 77-126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 18.7 | 93 | 77-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 22.0 | 110 | 80-138 | |
| 1,1-Dichloroethane | 20.0 | 18.9 | 94 | 76-120 | |
| 1,1-Dichloroethene | 20.0 | 18.5 | 92 | 75-135 | |
| 1,1-Dichloropropene | 20.0 | 21.0 | 105 | 80-120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 19.1 | 96 | 61-121 | |
| 1,2,3-Trichloropropane | 20.0 | 18.4 | 92 | 80-126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 19.0 | 95 | 64-124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 19.3 | 97 | 80-129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 18.5 | 92 | 61-132 | |
| 1,2-Dichlorobenzene | 20.0 | 18.8 | 94 | 76-120 | |
| 1,2-Dichloroethane | 20.0 | 20.7 | 104 | 72-120 | |
| 1,2-Dichloropropane | 20.0 | 17.5 | 87 | 80-120 | |
| 1,3-Dichlorobenzene | 20.0 | 18.4 | 92 | 78-120 | |
| 1,3-Dichloropropane | 20.0 | 18.4 | 92 | 80-120 | |
| 1,4-Dichlorobenzene | 20.0 | 18.1 | 91 | 75-120 | |
| 2,2-Dichloropropane | 20.0 | 23.0 | 115 | 69-135 | |
| 2-Butanone (MEK) | 40.0 | 30.7 | 77 | 52-131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 13.8 J | 69 | 28-173 | |
| 2-Chlorotoluene | 20.0 | 19.4 | 97 | 78-120 | |
| 2-Hexanone | 40.0 | 32.4 | 81 | 64-136 | |
| 4-Chlorotoluene | 20.0 | 19.4 | 97 | 79-120 | |
| p-Isopropyltoluene | 20.0 | 19.9 | 99 | 77-131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 31.6 | 79 | 67-135 | |
| Acetone | 40.0 | 33.8 | 84 | 41-137 | |
| Benzene | 20.0 | 18.6 | 93 | 79-120 | |
| Bromobenzene | 20.0 | 20.2 | 101 | 80-120 | |
| Bromoform | 20.0 | 17.5 | 88 | 62-133 | |
| Bromomethane | 20.0 | 18.6 | 93 | 42-136 | |
| Carbon disulfide | 20.0 | 16.2 | 81 | 62-146 | |
| Carbon tetrachloride | 20.0 | 22.9 | 115 | 71-129 | |
| Chlorobenzene | 20.0 | 19.5 | 98 | 78-120 | |
| Dibromochloromethane | 20.0 | 18.4 | 92 | 72-127 | |
| Chloroethane | 20.0 | 18.5 | 92 | 58-120 | |
| Chloroform | 20.0 | 20.8 | 104 | 77-120 | |
| Chloromethane | 20.0 | 14.3 | 72 | 50-120 | |
| cis-1,2-Dichloroethene | 20.0 | 20.5 | 102 | 76-120 | |
| cis-1,3-Dichloropropene | 20.0 | 17.9 | 89 | 74-128 | |
| Dibromomethane | 20.0 | 21.3 | 107 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146567.D

Lab ID: LCS 240-207560/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 20.0 | 20.9 | 105 | 79-120 | |
| Dichlorodifluoromethane | 20.0 | 15.4 | 77 | 26-120 | |
| Bromodichloromethane | 20.0 | 19.3 | 96 | 80-122 | |
| tert-Butyl alcohol (TBA) | 200 | 159 J | 80 | 60-125 | |
| Ethylbenzene | 20.0 | 20.0 | 100 | 79-120 | |
| Hexachlorobutadiene | 20.0 | 20.9 | 105 | 54-131 | |
| Isopropylbenzene | 20.0 | 20.5 | 102 | 76-122 | |
| Methyl tert-butyl ether | 20.0 | 18.8 | 94 | 49-165 | |
| Methylene Chloride | 20.0 | 15.3 | 76 | 75-120 | |
| m-Xylene & p-Xylene | 20.0 | 19.9 | 99 | 80-120 | |
| Naphthalene | 20.0 | 17.7 | 88 | 65-123 | |
| n-Butylbenzene | 20.0 | 19.4 | 97 | 68-135 | |
| N-Propylbenzene | 20.0 | 19.3 | 96 | 80-129 | |
| o-Xylene | 20.0 | 20.3 | 102 | 80-120 | |
| sec-Butylbenzene | 20.0 | 19.4 | 97 | 74-129 | |
| Styrene | 20.0 | 18.4 | 92 | 80-120 | |
| Tetrachloroethene | 20.0 | 20.7 | 104 | 79-120 | |
| Toluene | 20.0 | 19.9 | 100 | 75-120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.8 | 104 | 78-120 | |
| trans-1,3-Dichloropropene | 20.0 | 19.3 | 97 | 73-131 | |
| Trichloroethene | 20.0 | 20.4 | 102 | 79-120 | |
| Trichlorofluoromethane | 20.0 | 21.9 | 109 | 57-146 | |
| Vinyl acetate | 20.0 | 17.2 | 86 | 70-130 | |
| Vinyl chloride | 20.0 | 14.8 | 74 | 57-120 | |
| Xylenes, Total | 40.0 | 40.2 | 101 | 80-120 | |
| tert-Butylbenzene | 20.0 | 19.2 | 96 | 76-126 | |
| 1,2-Dibromoethane | 20.0 | 20.9 | 104 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: UX80877.D

Lab ID: LCS 240-207720/6 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 50.0 | 49.7 | 99 | 80-120 | |
| 1,1,1-Trichloroethane | 50.0 | 49.6 | 99 | 77-126 | |
| 1,1,2,2-Tetrachloroethane | 50.0 | 49.8 | 100 | 77-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 50.0 | 51.8 | 104 | 80-138 | |
| 1,1-Dichloroethane | 50.0 | 48.9 | 98 | 76-120 | |
| 1,1-Dichloroethene | 50.0 | 48.3 | 97 | 75-135 | |
| 1,1-Dichloropropene | 50.0 | 52.2 | 104 | 80-120 | |
| 1,2,3-Trichlorobenzene | 50.0 | 45.0 | 90 | 61-121 | |
| 1,2,3-Trichloropropane | 50.0 | 52.8 | 106 | 80-126 | |
| 1,2,4-Trichlorobenzene | 50.0 | 46.3 | 93 | 64-124 | |
| 1,2,4-Trimethylbenzene | 50.0 | 49.4 | 99 | 80-129 | |
| 1,2-Dibromo-3-Chloropropane | 50.0 | 48.6 | 97 | 61-132 | |
| 1,2-Dichlorobenzene | 50.0 | 46.6 | 93 | 76-120 | |
| 1,2-Dichloroethane | 50.0 | 49.9 | 100 | 72-120 | |
| 1,2-Dichloropropane | 50.0 | 47.1 | 94 | 80-120 | |
| 1,3-Dichlorobenzene | 50.0 | 45.9 | 92 | 78-120 | |
| 1,3-Dichloropropane | 50.0 | 46.0 | 92 | 80-120 | |
| 1,4-Dichlorobenzene | 50.0 | 45.5 | 91 | 75-120 | |
| 2,2-Dichloropropane | 50.0 | 44.2 | 88 | 69-135 | |
| 2-Butanone (MEK) | 100 | 98.6 | 99 | 52-131 | |
| 2-Chloroethyl vinyl ether | 50.0 | NC | NaN | 28-173 | |
| 2-Chlorotoluene | 50.0 | 48.0 | 96 | 78-120 | |
| 2-Hexanone | 100 | 102 | 102 | 64-136 | |
| 4-Chlorotoluene | 50.0 | 45.9 | 92 | 79-120 | |
| p-Isopropyltoluene | 50.0 | 50.3 | 101 | 77-131 | |
| 4-Methyl-2-pentanone (MIBK) | 100 | 107 | 107 | 67-135 | |
| Acetone | 100 | 111 | 111 | 41-137 | |
| Benzene | 50.0 | 46.8 | 94 | 79-120 | |
| Bromobenzene | 50.0 | 46.6 | 93 | 80-120 | |
| Bromoform | 50.0 | 49.8 | 100 | 62-133 | |
| Bromomethane | 50.0 | 56.1 | 112 | 42-136 | |
| Carbon disulfide | 50.0 | 50.4 | 101 | 62-146 | |
| Carbon tetrachloride | 50.0 | 53.3 | 107 | 71-129 | |
| Chlorobenzene | 50.0 | 45.5 | 91 | 78-120 | |
| Dibromochloromethane | 50.0 | 52.7 | 105 | 72-127 | |
| Chloroethane | 50.0 | 49.4 | 99 | 58-120 | |
| Chloroform | 50.0 | 49.5 | 99 | 77-120 | |
| Chloromethane | 50.0 | 50.8 | 102 | 50-120 | |
| cis-1,2-Dichloroethene | 50.0 | 48.3 | 97 | 76-120 | |
| cis-1,3-Dichloropropene | 50.0 | 48.3 | 97 | 74-128 | |
| Dibromomethane | 50.0 | 49.9 | 100 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: UX80877.D

Lab ID: LCS 240-207720/6 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 50.0 | 49.5 | 99 | 79-120 | |
| Dichlorodifluoromethane | 50.0 | 52.4 | 105 | 26-120 | |
| Bromodichloromethane | 50.0 | 50.4 | 101 | 80-122 | |
| tert-Butyl alcohol (TBA) | 500 | 440 | 88 | 60-125 | |
| Ethylbenzene | 50.0 | 45.5 | 91 | 79-120 | |
| Hexachlorobutadiene | 50.0 | 48.6 | 97 | 54-131 | |
| Isopropylbenzene | 50.0 | 50.1 | 100 | 76-122 | |
| Methyl tert-butyl ether | 50.0 | 47.8 | 96 | 49-165 | |
| Methylene Chloride | 50.0 | 52.3 | 105 | 75-120 | |
| m-Xylene & p-Xylene | 50.0 | 46.2 | 92 | 80-120 | |
| Naphthalene | 50.0 | 52.0 | 104 | 65-123 | |
| n-Butylbenzene | 50.0 | 50.0 | 100 | 68-135 | |
| N-Propylbenzene | 50.0 | 48.6 | 97 | 80-129 | |
| o-Xylene | 50.0 | 47.8 | 96 | 80-120 | |
| sec-Butylbenzene | 50.0 | 51.0 | 102 | 74-129 | |
| Styrene | 50.0 | 45.9 | 92 | 80-120 | |
| Tetrachloroethene | 50.0 | 45.4 | 91 | 79-120 | |
| Toluene | 50.0 | 45.5 | 91 | 75-120 | |
| trans-1,2-Dichloroethene | 50.0 | 48.7 | 97 | 78-120 | |
| trans-1,3-Dichloropropene | 50.0 | 42.9 | 86 | 73-131 | |
| Trichloroethene | 50.0 | 48.9 | 98 | 79-120 | |
| Trichlorofluoromethane | 50.0 | 50.8 | 102 | 57-146 | |
| Vinyl acetate | 50.0 | 39.2 | 78 | 70-130 | |
| Vinyl chloride | 50.0 | 49.9 | 100 | 57-120 | |
| Xylenes, Total | 100 | 94.0 | 94 | 80-120 | |
| tert-Butylbenzene | 50.0 | 55.3 | 111 | 76-126 | |
| 1,2-Dibromoethane | 50.0 | 48.6 | 97 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXX3079.D

Lab ID: 200-30727-B-1 MS Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 500 | 14 U | 486 | 97 | 72-120 | |
| 1,1,1-Trichloroethane | 500 | 22 U | 466 | 93 | 69-122 | |
| 1,1,2,2-Tetrachloroethane | 500 | 11 U | 484 | 97 | 61-130 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 500 | 23 U | 502 | 100 | 44-140 | |
| 1,1-Dichloroethane | 500 | 15 U | 477 | 95 | 73-124 | |
| 1,1-Dichloroethene | 500 | 23 U | 485 | 97 | 67-124 | |
| 1,1-Dichloropropene | 500 | 21 U | 526 | 105 | 81-126 | |
| 1,2,3-Trichlorobenzene | 500 | 19 U | 409 | 82 | 46-120 | |
| 1,2,3-Trichloropropane | 500 | 22 U | 498 | 100 | 58-135 | |
| 1,2,4-Trichlorobenzene | 500 | 16 U | 426 | 85 | 48-120 | |
| 1,2,4-Trimethylbenzene | 500 | 21 U | 477 | 95 | 64-122 | |
| 1,2-Dibromo-3-Chloropropane | 500 | 41 U | 381 | 76 | 42-130 | |
| 1,2-Dichlorobenzene | 500 | 13 U | 475 | 95 | 67-118 | |
| 1,2-Dichloroethane | 500 | 12 U | 489 | 98 | 74-125 | |
| 1,2-Dichloropropane | 500 | 13 U | 483 | 97 | 73-122 | |
| 1,3-Dichlorobenzene | 500 | 9.5 U | 468 | 94 | 65-120 | |
| 1,3-Dichloropropane | 500 | 9.5 U | 483 | 97 | 59-133 | |
| 1,4-Dichlorobenzene | 500 | 14 U | 472 | 94 | 66-120 | |
| 2,2-Dichloropropane | 500 | 13 U | 402 | 80 | 48-126 | |
| 2-Butanone (MEK) | 1000 | 27 U | 953 | 95 | 49-132 | |
| 2-Chloroethyl vinyl ether | 500 | 30 U | 278 J | 56 | 10-150 | |
| 2-Chlorotoluene | 500 | 20 U | 484 | 97 | 70-122 | |
| 2-Hexanone | 1000 | 24 U | 954 | 95 | 49-142 | |
| 4-Chlorotoluene | 500 | 15 U | 501 | 100 | 74-120 | |
| p-Isopropyltoluene | 500 | 22 U | 475 | 95 | 61-123 | |
| 4-Methyl-2-pentanone (MIBK) | 1000 | 50 U | 957 | 96 | 58-136 | |
| Acetone | 1000 | 47 U | 998 | 100 | 32-126 | |
| Benzene | 500 | 18 U | 471 | 94 | 73-121 | |
| Bromobenzene | 500 | 18 U | 486 | 97 | 73-120 | |
| Bromoform | 500 | 28 U | 419 | 84 | 45-121 | |
| Bromomethane | 500 | 22 U | 402 | 80 | 26-136 | |
| Carbon disulfide | 500 | 19 U | 465 | 93 | 54-144 | |
| Carbon tetrachloride | 500 | 22 U | 475 | 95 | 65-129 | |
| Chlorobenzene | 500 | 13 U | 482 | 96 | 72-120 | |
| Dibromochloromethane | 500 | 22 U | 458 | 92 | 62-122 | |
| Chloroethane | 500 | 16 U | 426 | 85 | 27-131 | |
| Chloroform | 500 | 13 U | 488 | 98 | 73-121 | |
| Chloromethane | 500 | 22 U | 356 | 71 | 39-134 | |
| cis-1,2-Dichloroethene | 500 | 1200 | 1660 | 95 | 66-124 | |
| cis-1,3-Dichloropropene | 500 | 23 U | 470 | 94 | 60-120 | |
| Dibromomethane | 500 | 21 U | 479 | 96 | 63-134 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXX3079.D

Lab ID: 200-30727-B-1 MS Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Bromochloromethane | 500 | 25 U | 481 | 96 | 60-132 | |
| Dichlorodifluoromethane | 500 | 16 U | 360 | 72 | 14-137 | |
| Bromodichloromethane | 500 | 15 U | 458 | 92 | 72-120 | |
| Ethylbenzene | 500 | 13 U | 489 | 98 | 68-121 | |
| Hexachlorobutadiene | 500 | 18 U | 395 | 79 | 48-120 | |
| Isopropylbenzene | 500 | 18 U | 498 | 100 | 61-122 | |
| Methyl tert-butyl ether | 500 | 10 U | 443 | 89 | 61-121 | |
| Methylene Chloride | 500 | 52 | 578 | 105 | 70-124 | |
| m-Xylene & p-Xylene | 500 | 12 U | 509 | 102 | 66-123 | |
| Naphthalene | 500 | 23 U | 498 | 100 | 38-128 | |
| n-Butylbenzene | 500 | 16 U | 473 | 95 | 54-126 | |
| N-Propylbenzene | 500 | 20 U | 516 | 103 | 63-129 | |
| o-Xylene | 500 | 13 U | 501 | 100 | 68-121 | |
| sec-Butylbenzene | 500 | 24 U | 489 | 98 | 61-123 | |
| Styrene | 500 | 23 U | 487 | 97 | 64-126 | |
| Tetrachloroethene | 500 | 660 | 1140 | 96 | 59-125 | |
| Toluene | 500 | 12 U | 483 | 97 | 72-122 | |
| trans-1,2-Dichloroethene | 500 | 15 U | 493 | 99 | 72-125 | |
| trans-1,3-Dichloropropene | 500 | 28 U | 474 | 95 | 58-132 | |
| Trichloroethene | 500 | 590 | 1110 | 103 | 61-129 | |
| Trichlorofluoromethane | 500 | 25 U | 455 | 91 | 49-133 | |
| Vinyl acetate | 500 | 21 U | 376 | 75 | 26-189 | |
| Vinyl chloride | 500 | 290 | 693 | 80 | 44-122 | |
| Xylenes, Total | 1000 | 26 U | 1010 | 101 | 67-122 | |
| tert-Butylbenzene | 500 | 21 U | 484 | 97 | 63-122 | |
| 1,2-Dibromoethane | 500 | 16 U | 494 | 99 | 69-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146590.D

Lab ID: 240-57921-A-4-J MS Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 20.2 | 0.65 U | 20.0 | 99 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 20.2 | 0.31 U | 0.31 U | 0 | 16-179 | F1 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.2 | 0.42 U | 21.1 | 104 | 50-147 | |
| 1,1-Dichloroethane | 20.2 | 0.34 U | 16.3 | 80 | 54-122 | |
| 1,1-Dichloroethene | 20.2 | 0.79 U | 24.7 | 122 | 49-157 | |
| 1,2,4-Trichlorobenzene | 20.2 | 0.38 U | 9.56 | 47 | 10-120 | |
| 1,2-Dibromo-3-Chloropropane | 20.2 | 2.7 U | 6.61 J | 33 | 10-153 | |
| 1,2-Dichlorobenzene | 20.2 | 0.34 U | 11.6 | 57 | 17-122 | |
| 1,2-Dichloroethane | 20.2 | 0.46 U | 18.3 | 90 | 49-123 | |
| 1,2-Dichloropropane | 20.2 | 0.14 U | 15.3 | 75 | 61-120 | |
| 1,3-Dichlorobenzene | 20.2 | 0.69 U | 10.5 | 52 | 16-126 | |
| 1,4-Dichlorobenzene | 20.2 | 0.28 U | 10.3 | 51 | 15-121 | |
| 2-Butanone (MEK) | 40.5 | 2.9 J | 32.2 | 72 | 30-143 | |
| 2-Hexanone | 40.5 | 1.3 J | 29.9 | 71 | 37-147 | |
| 4-Methyl-2-pentanone (MIBK) | 40.5 | 1.2 U | 28.0 | 69 | 43-147 | |
| Acetone | 40.5 | 34 | 78.3 | 110 | 24-140 | |
| Benzene | 20.2 | 0.89 U | 16.0 | 79 | 53-120 | |
| Bromoform | 20.2 | 0.24 U | 14.0 | 69 | 18-129 | |
| Bromomethane | 20.2 | 0.42 U | 12.4 | 61 | 33-130 | |
| Carbon disulfide | 20.2 | 2.2 J | 15.1 | 64 | 20-151 | |
| Carbon tetrachloride | 20.2 | 0.65 U | 19.5 | 96 | 32-137 | |
| Chlorobenzene | 20.2 | 0.53 U | 14.1 | 69 | 37-120 | |
| Dibromochloromethane | 20.2 | 0.37 U | 14.4 | 71 | 29-135 | |
| Chloroethane | 20.2 | 0.44 U | 15.8 | 78 | 45-120 | |
| Chloroform | 20.2 | 0.36 U | 18.3 | 90 | 53-120 | |
| Chloromethane | 20.2 | 0.78 U | 12.2 | 60 | 34-120 | |
| cis-1,2-Dichloroethene | 20.2 | 0.28 U | 17.3 | 85 | 50-120 | |
| cis-1,3-Dichloropropene | 20.2 | 0.71 U | 9.86 | 49 | 27-133 | |
| Dichlorodifluoromethane | 20.2 | 0.28 U | 13.8 | 68 | 17-120 | |
| Bromodichloromethane | 20.2 | 0.29 U | 14.2 | 70 | 35-132 | |
| Ethylbenzene | 20.2 | 0.32 U | 14.0 | 69 | 30-131 | |
| Isopropylbenzene | 20.2 | 0.20 U | 13.3 | 66 | 21-134 | |
| Methyl tert-butyl ether | 20.2 | 0.43 U | 17.4 | 86 | 51-157 | |
| Methylene Chloride | 20.2 | 0.73 U | 16.2 | 80 | 54-120 | |
| m-Xylene & p-Xylene | 20.2 | 0.42 U | 13.6 | 67 | 29-131 | |
| o-Xylene | 20.2 | 0.57 U | 14.4 | 71 | 29-134 | |
| Styrene | 20.2 | 0.40 U | 12.7 | 63 | 27-127 | |
| Tetrachloroethene | 20.2 | 0.79 U | 15.3 | 76 | 31-135 | |
| Toluene | 20.2 | 0.40 J | 15.9 | 76 | 39-129 | |
| trans-1,2-Dichloroethene | 20.2 | 0.32 U | 17.3 | 85 | 50-123 | |
| trans-1,3-Dichloropropene | 20.2 | 0.29 U | 12.3 | 61 | 28-137 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146590.D

Lab ID: 240-57921-A-4-J MS Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Trichloroethene | 20.2 | 0.38 U | 31.6 | 156 | 10-177 | |
| Trichlorofluoromethane | 20.2 | 0.29 U | 18.7 | 93 | 36-142 | |
| Vinyl chloride | 20.2 | 0.30 U | 13.6 | 67 | 42-120 | |
| Xylenes, Total | 40.5 | 0.54 U | 28.0 | 69 | 30-131 | |
| 1,2-Dibromoethane | 20.2 | 0.53 U | 17.1 | 85 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: UX80889.D

Lab ID: 240-57972-C-3-C MS Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Acetone | 54.2 | 6.4 U | 37.4 | 69 | 24-140 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXX3080.D

Lab ID: 200-30727-B-1 MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 500 | 467 | 93 | 4 | 22 | 72-120 | |
| 1,1,1-Trichloroethane | 500 | 479 | 96 | 3 | 14 | 69-122 | |
| 1,1,2,2-Tetrachloroethane | 500 | 485 | 97 | 0 | 18 | 61-130 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 500 | 518 | 104 | 3 | 35 | 44-140 | |
| 1,1-Dichloroethane | 500 | 479 | 96 | 0 | 14 | 73-124 | |
| 1,1-Dichloroethene | 500 | 504 | 101 | 4 | 24 | 67-124 | |
| 1,1-Dichloropropene | 500 | 514 | 103 | 2 | 11 | 81-126 | |
| 1,2,3-Trichlorobenzene | 500 | 414 | 83 | 1 | 35 | 46-120 | |
| 1,2,3-Trichloropropane | 500 | 491 | 98 | 1 | 25 | 58-135 | |
| 1,2,4-Trichlorobenzene | 500 | 425 | 85 | 0 | 28 | 48-120 | |
| 1,2,4-Trimethylbenzene | 500 | 470 | 94 | 2 | 14 | 64-122 | |
| 1,2-Dibromo-3-Chloropropane | 500 | 387 | 77 | 2 | 24 | 42-130 | |
| 1,2-Dichlorobenzene | 500 | 465 | 93 | 2 | 15 | 67-118 | |
| 1,2-Dichloroethane | 500 | 487 | 97 | 0 | 24 | 74-125 | |
| 1,2-Dichloropropane | 500 | 486 | 97 | 1 | 15 | 73-122 | |
| 1,3-Dichlorobenzene | 500 | 469 | 94 | 0 | 15 | 65-120 | |
| 1,3-Dichloropropane | 500 | 483 | 97 | 0 | 35 | 59-133 | |
| 1,4-Dichlorobenzene | 500 | 465 | 93 | 1 | 16 | 66-120 | |
| 2,2-Dichloropropane | 500 | 410 | 82 | 2 | 30 | 48-126 | |
| 2-Butanone (MEK) | 1000 | 997 | 100 | 4 | 19 | 49-132 | |
| 2-Chloroethyl vinyl ether | 500 | 254 J | 51 | 9 | 30 | 10-150 | |
| 2-Chlorotoluene | 500 | 485 | 97 | 0 | 18 | 70-122 | |
| 2-Hexanone | 1000 | 985 | 98 | 3 | 27 | 49-142 | |
| 4-Chlorotoluene | 500 | 486 | 97 | 3 | 15 | 74-120 | |
| p-Isopropyltoluene | 500 | 477 | 95 | 0 | 27 | 61-123 | |
| 4-Methyl-2-pentanone (MIBK) | 1000 | 965 | 96 | 1 | 32 | 58-136 | |
| Acetone | 1000 | 1070 | 107 | 7 | 28 | 32-126 | |
| Benzene | 500 | 469 | 94 | 1 | 13 | 73-121 | |
| Bromobenzene | 500 | 494 | 99 | 2 | 13 | 73-120 | |
| Bromoform | 500 | 416 | 83 | 1 | 19 | 45-121 | |
| Bromomethane | 500 | 420 | 84 | 4 | 35 | 26-136 | |
| Carbon disulfide | 500 | 455 | 91 | 2 | 34 | 54-144 | |
| Carbon tetrachloride | 500 | 485 | 97 | 2 | 20 | 65-129 | |
| Chlorobenzene | 500 | 477 | 95 | 1 | 15 | 72-120 | |
| Dibromochloromethane | 500 | 464 | 93 | 1 | 19 | 62-122 | |
| Chloroethane | 500 | 432 | 86 | 1 | 35 | 27-131 | |
| Chloroform | 500 | 492 | 98 | 1 | 17 | 73-121 | |
| Chloromethane | 500 | 379 | 76 | 6 | 20 | 39-134 | |
| cis-1,2-Dichloroethene | 500 | 1660 | 94 | 0 | 22 | 66-124 | |
| cis-1,3-Dichloropropene | 500 | 483 | 97 | 3 | 21 | 60-120 | |
| Dibromomethane | 500 | 478 | 96 | 0 | 35 | 63-134 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXX3080.D

Lab ID: 200-30727-B-1 MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Bromochloromethane | 500 | 467 | 93 | 3 | 35 | 60-132 | |
| Dichlorodifluoromethane | 500 | 354 | 71 | 2 | 34 | 14-137 | |
| Bromodichloromethane | 500 | 462 | 92 | 1 | 19 | 72-120 | |
| Ethylbenzene | 500 | 489 | 98 | 0 | 16 | 68-121 | |
| Hexachlorobutadiene | 500 | 410 | 82 | 4 | 34 | 48-120 | |
| Isopropylbenzene | 500 | 479 | 96 | 4 | 20 | 61-122 | |
| Methyl tert-butyl ether | 500 | 457 | 91 | 3 | 12 | 61-121 | |
| Methylene Chloride | 500 | 579 | 106 | 0 | 14 | 70-124 | |
| m-Xylene & p-Xylene | 500 | 488 | 98 | 4 | 15 | 66-123 | |
| Naphthalene | 500 | 431 | 86 | 14 | 35 | 38-128 | |
| n-Butylbenzene | 500 | 463 | 93 | 2 | 33 | 54-126 | |
| N-Propylbenzene | 500 | 510 | 102 | 1 | 29 | 63-129 | |
| o-Xylene | 500 | 492 | 98 | 2 | 14 | 68-121 | |
| sec-Butylbenzene | 500 | 480 | 96 | 2 | 29 | 61-123 | |
| Styrene | 500 | 464 | 93 | 5 | 15 | 64-126 | |
| Tetrachloroethene | 500 | 1160 | 100 | 2 | 20 | 59-125 | |
| Toluene | 500 | 478 | 96 | 1 | 15 | 72-122 | |
| trans-1,2-Dichloroethene | 500 | 486 | 97 | 1 | 25 | 72-125 | |
| trans-1,3-Dichloropropene | 500 | 494 | 99 | 4 | 22 | 58-132 | |
| Trichloroethene | 500 | 1110 | 103 | 0 | 14 | 61-129 | |
| Trichlorofluoromethane | 500 | 479 | 96 | 5 | 25 | 49-133 | |
| Vinyl acetate | 500 | 383 | 77 | 2 | 15 | 26-189 | |
| Vinyl chloride | 500 | 716 | 84 | 3 | 35 | 44-122 | |
| Xylenes, Total | 1000 | 980 | 98 | 3 | 14 | 67-122 | |
| tert-Butylbenzene | 500 | 485 | 97 | 0 | 24 | 63-122 | |
| 1,2-Dibromoethane | 500 | 485 | 97 | 2 | 24 | 69-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146591.D

Lab ID: 240-57921-A-4-K MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 19.9 | 20.2 | 101 | 1 | 30 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 19.9 | 0.31 U | 0 | NC | 30 | 16-179 | F1 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 19.9 | 17.4 | 87 | 19 | 30 | 50-147 | |
| 1,1-Dichloroethane | 19.9 | 16.3 | 82 | 0 | 30 | 54-122 | |
| 1,1-Dichloroethene | 19.9 | 22.7 | 114 | 8 | 30 | 49-157 | |
| 1,2,4-Trichlorobenzene | 19.9 | 10.9 | 55 | 13 | 30 | 10-120 | |
| 1,2-Dibromo-3-Chloropropane | 19.9 | 5.77 J | 29 | 14 | 30 | 10-153 | |
| 1,2-Dichlorobenzene | 19.9 | 13.2 | 66 | 13 | 30 | 17-122 | |
| 1,2-Dichloroethane | 19.9 | 17.9 | 90 | 2 | 30 | 49-123 | |
| 1,2-Dichloropropane | 19.9 | 15.8 | 79 | 4 | 30 | 61-120 | |
| 1,3-Dichlorobenzene | 19.9 | 12.7 | 64 | 19 | 30 | 16-126 | |
| 1,4-Dichlorobenzene | 19.9 | 11.8 | 59 | 13 | 30 | 15-121 | |
| 2-Butanone (MEK) | 39.9 | 28.6 | 64 | 12 | 30 | 30-143 | |
| 2-Hexanone | 39.9 | 28.2 | 67 | 6 | 30 | 37-147 | |
| 4-Methyl-2-pentanone (MIBK) | 39.9 | 27.1 | 68 | 3 | 30 | 43-147 | |
| Acetone | 39.9 | 71.4 | 94 | 9 | 30 | 24-140 | |
| Benzene | 19.9 | 16.5 | 83 | 3 | 30 | 53-120 | |
| Bromoform | 19.9 | 13.5 | 68 | 4 | 30 | 18-129 | |
| Bromomethane | 19.9 | 11.7 | 59 | 6 | 30 | 33-130 | |
| Carbon disulfide | 19.9 | 15.5 | 67 | 2 | 30 | 20-151 | |
| Carbon tetrachloride | 19.9 | 20.0 | 100 | 2 | 30 | 32-137 | |
| Chlorobenzene | 19.9 | 14.9 | 75 | 6 | 30 | 37-120 | |
| Dibromochloromethane | 19.9 | 13.5 | 68 | 6 | 30 | 29-135 | |
| Chloroethane | 19.9 | 16.7 | 84 | 5 | 30 | 45-120 | |
| Chloroform | 19.9 | 18.6 | 93 | 2 | 30 | 53-120 | |
| Chloromethane | 19.9 | 12.1 | 61 | 0 | 30 | 34-120 | |
| cis-1,2-Dichloroethene | 19.9 | 17.6 | 88 | 2 | 30 | 50-120 | |
| cis-1,3-Dichloropropene | 19.9 | 9.38 | 47 | 5 | 30 | 27-133 | |
| Dichlorodifluoromethane | 19.9 | 13.7 | 69 | 1 | 30 | 17-120 | |
| Bromodichloromethane | 19.9 | 13.6 | 68 | 4 | 30 | 35-132 | |
| Ethylbenzene | 19.9 | 15.4 | 77 | 9 | 30 | 30-131 | |
| Isopropylbenzene | 19.9 | 15.4 | 77 | 14 | 30 | 21-134 | |
| Methyl tert-butyl ether | 19.9 | 17.2 | 86 | 2 | 30 | 51-157 | |
| Methylene Chloride | 19.9 | 17.4 | 87 | 7 | 30 | 54-120 | |
| m-Xylene & p-Xylene | 19.9 | 14.9 | 75 | 9 | 30 | 29-131 | |
| o-Xylene | 19.9 | 15.7 | 79 | 9 | 30 | 29-134 | |
| Styrene | 19.9 | 13.4 | 67 | 5 | 30 | 27-127 | |
| Tetrachloroethene | 19.9 | 16.8 | 84 | 9 | 30 | 31-135 | |
| Toluene | 19.9 | 16.7 | 82 | 5 | 30 | 39-129 | |
| trans-1,2-Dichloroethene | 19.9 | 17.6 | 88 | 2 | 30 | 50-123 | |
| trans-1,3-Dichloropropene | 19.9 | 11.4 | 57 | 8 | 30 | 28-137 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146591.D

Lab ID: 240-57921-A-4-K MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Trichloroethene | 19.9 | 32.5 | 163 | 3 | 30 | 10-177 | |
| Trichlorofluoromethane | 19.9 | 19.5 | 98 | 4 | 30 | 36-142 | |
| Vinyl chloride | 19.9 | 13.7 | 69 | 1 | 30 | 42-120 | |
| Xylenes, Total | 39.9 | 30.6 | 77 | 9 | 30 | 30-131 | |
| 1,2-Dibromoethane | 19.9 | 16.6 | 83 | 3 | 30 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: UX80890.D

Lab ID: 240-57972-C-3-D MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Acetone | 56.7 | 46.4 | 82 | 21 | 30 | 24-140 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab File ID: 146535.D Lab Sample ID: MB 240-207343/6

Matrix: Solid Heated Purge:(Y/N) Y

Instrument ID: A3UX14 Date Analyzed: 11/18/2015 12:20

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-207343/5 | 146534.D | 11/18/2015 11:58 |
| | LCS 240-207343/7 | 146536.D | 11/18/2015 12:41 |
| F-VS-UST-P-1 | 240-57930-1 | 146551.D | 11/18/2015 18:03 |
| F-VS-UST-P-3 | 240-57930-3 | 146552.D | 11/18/2015 18:25 |
| F-VB-UST-P-1 | 240-57930-5 | 146556.D | 11/18/2015 19:51 |
| F-VB-UST-P-2 | 240-57930-6 | 146557.D | 11/18/2015 20:12 |
| F-VB-UST-P-3 | 240-57930-7 | 146558.D | 11/18/2015 20:33 |
| F-VB-UST-P-4 | 240-57930-8 | 146559.D | 11/18/2015 20:55 |
| F-VB-UST-P-5 | 240-57930-9 | 146560.D | 11/18/2015 21:16 |
| F-VB-UST-P-6 | 240-57930-10 | 146561.D | 11/18/2015 21:38 |
| F-VS-UST-P-DUP | 240-57930-11 | 146562.D | 11/18/2015 21:59 |

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab File ID: 146568.D Lab Sample ID: MB 240-207560/6

Matrix: Solid Heated Purge:(Y/N) Y

Instrument ID: A3UX14 Date Analyzed: 11/19/2015 12:29

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-207560/5 | 146567.D | 11/19/2015 12:08 |
| | LCS 240-207560/7 | 146569.D | 11/19/2015 12:51 |
| F-VS-UST-P-4 | 240-57930-4 | 146581.D | 11/19/2015 17:08 |
| F-VS-UST-P-2 | 240-57930-2 | 146583.D | 11/19/2015 17:52 |
| F-VS-UST-P-2 | 240-57930-2 | 146584.D | 11/19/2015 18:13 |
| | 240-57921-A-4-J MS | 146590.D | 11/19/2015 20:22 |
| | 240-57921-A-4-K MSD | 146591.D | 11/19/2015 20:43 |

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab File ID: UX80879.D Lab Sample ID: MB 240-207720/8

Matrix: Solid Heated Purge:(Y/N) Y

Instrument ID: A3UX8 Date Analyzed: 11/20/2015 05:08

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-207720/6 | UX80877.D | 11/20/2015 04:26 |
| | 240-57972-C-3-C MS | UX80889.D | 11/20/2015 08:41 |
| | 240-57972-C-3-D MSD | UX80890.D | 11/20/2015 09:02 |
| F-VS-UST-P-2 | 240-57930-2 | UX80893.D | 11/20/2015 10:06 |

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: UXX3070.D  Lab Sample ID: MB 240-207096/6

Matrix: Water  Heated Purge:(Y/N) N

Instrument ID: A3UX10  Date Analyzed: 11/17/2015 12:17

GC Column: DB-624  ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-207096/4 | UXX3068.D | 11/17/2015 11:34 |
| TB-111315 | 240-57930-12 | UXX3077.D | 11/17/2015 14:51 |
| | 200-30727-B-1 MS | UXX3079.D | 11/17/2015 15:35 |
| | 200-30727-B-1 MSD | UXX3080.D | 11/17/2015 15:57 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB7317.D  BFB Injection Date: 11/16/2015

Instrument ID: A3UX10  BFB Injection Time: 09:00

Analysis Batch No.: 206893

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 20.7 | |
| 75 | 30.0 - 60.0 % of mass 95 | 52.0 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.9 | |
| 173 | Less than 2.0 % of mass 174 | 0.0 | (0.0) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 100.1 | |
| 175 | 5.0 - 9.0 % of mass 174 | 7.3 | (7.3) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 95.9 | (95.7) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 6.6 | (6.9) 2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD8260 240-206893/2 | UXX3039.D | 11/16/2015 | 09:23 |
| | STD8260 240-206893/3 | UXX3040.D | 11/16/2015 | 09:45 |
| | STD8260 240-206893/4 | UXX3041.D | 11/16/2015 | 10:07 |
| | STD8260 240-206893/5 | UXX3042.D | 11/16/2015 | 10:29 |
| | STD8260 240-206893/6 | UXX3043.D | 11/16/2015 | 10:51 |
| | STD8260 240-206893/7 | UXX3044.D | 11/16/2015 | 11:13 |
| | STD8260 240-206893/8 | UXX3045.D | 11/16/2015 | 11:35 |
| | STDA9 240-206893/10 | UXX3047.D | 11/16/2015 | 12:19 |
| | STDA9 240-206893/11 | UXX3048.D | 11/16/2015 | 12:43 |
| | STDA9 240-206893/12 | UXX3049.D | 11/16/2015 | 13:05 |
| | STDA9 240-206893/13 | UXX3050.D | 11/16/2015 | 13:27 |
| | STDA9 240-206893/14 | UXX3051.D | 11/16/2015 | 13:49 |
| | STDA9 240-206893/15 | UXX3052.D | 11/16/2015 | 14:10 |
| | ICV 240-206893/16 | UXX3053.D | 11/16/2015 | 14:33 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB7318.D  BFB Injection Date: 11/17/2015

Instrument ID: A3UX10  BFB Injection Time: 10:26

Analysis Batch No.: 207096

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 23.5 | |
| 75 | 30.0 - 60.0 % of mass 95 | 58.5 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.9 | |
| 173 | Less than 2.0 % of mass 174 | 0.3 | (0.3) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 96.7 | |
| 175 | 5.0 - 9.0 % of mass 174 | 7.3 | (7.6) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 96.7 | (100.1) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 6.1 | (6.3) 2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-207096/2 | UXX3066.D | 11/17/2015 | 10:50 |
| | CCV 240-207096/3 | UXX3067.D | 11/17/2015 | 11:11 |
| | LCS 240-207096/4 | UXX3068.D | 11/17/2015 | 11:34 |
| | MB 240-207096/6 | UXX3070.D | 11/17/2015 | 12:17 |
| TB-111315 | 240-57930-12 | UXX3077.D | 11/17/2015 | 14:51 |
| | 200-30727-B-1 MS | UXX3079.D | 11/17/2015 | 15:35 |
| | 200-30727-B-1 MSD | UXX3080.D | 11/17/2015 | 15:57 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB14045.D BFB Injection Date: 08/13/2015

Instrument ID: A3UX14 BFB Injection Time: 13:54

Analysis Batch No.: 193150

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 19.1 | |
| 75 | 30.0 - 60.0 % of mass 95 | 45.9 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.7 | |
| 173 | Less than 2.0 % of mass 174 | 0.0 | (0.1) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 70.4 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.1 | (7.3) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 67.7 | (96.1) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.5 | (6.6) 2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | IC 240-193150/12 | 144576.D | 08/13/2015 | 17:50 |
| | IC 240-193150/13 | 144577.D | 08/13/2015 | 18:11 |
| | IC 240-193150/14 | 144578.D | 08/13/2015 | 18:33 |
| | IC 240-193150/15 | 144579.D | 08/13/2015 | 18:54 |
| | IC 240-193150/16 | 144580.D | 08/13/2015 | 19:16 |
| | IC 240-193150/17 | 144581.D | 08/13/2015 | 19:37 |
| | IC 240-193150/18 | 144582.D | 08/13/2015 | 19:59 |
| | IC 240-193150/19 | 144583.D | 08/13/2015 | 20:20 |
| | IC 240-193150/20 | 144584.D | 08/13/2015 | 20:42 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB14070.D BFB Injection Date: 09/02/2015

Instrument ID: A3UX14 BFB Injection Time: 13:49

Analysis Batch No.: 196026

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 21.6 | |
| 75 | 30.0 - 60.0 % of mass 95 | 45.0 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 7.0 | |
| 173 | Less than 2.0 % of mass 174 | 0.6 | (0.8) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 70.3 | |
| 175 | 5.0 - 9.0 % of mass 174 | 4.9 | (7.0) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 68.0 | (96.8) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.6 | (6.7) 2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | IC 240-196026/3 | 144935.D | 09/02/2015 | 14:32 |
| | IC 240-196026/4 | 144936.D | 09/02/2015 | 14:54 |
| | IC 240-196026/5 | 144937.D | 09/02/2015 | 15:15 |
| | ICIS 240-196026/6 | 144938.D | 09/02/2015 | 15:37 |
| | IC 240-196026/7 | 144939.D | 09/02/2015 | 15:58 |
| | IC 240-196026/8 | 144940.D | 09/02/2015 | 16:20 |
| | IC 240-196026/9 | 144941.D | 09/02/2015 | 16:41 |
| | IC 240-196026/10 | 144942.D | 09/02/2015 | 17:03 |
| | IC 240-196026/11 | 144943.D | 09/02/2015 | 17:24 |
| | ICV 240-196026/13 | 144945.D | 09/02/2015 | 18:07 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB14134.D BFB Injection Date: 11/18/2015

Instrument ID: A3UX14 BFB Injection Time: 10:30

Analysis Batch No.: 207343

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 19.7 | |
| 75 | 30.0 - 60.0 % of mass 95 | 46.5 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.7 | |
| 173 | Less than 2.0 % of mass 174 | 0.2 | (0.3) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 76.8 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.9 | (7.7) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 74.8 | (97.4) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.6 | (6.1) 2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-207343/3 | 146532.D | 11/18/2015 | 11:15 |
| | CCV 240-207343/4 | 146533.D | 11/18/2015 | 11:37 |
| | LCS 240-207343/5 | 146534.D | 11/18/2015 | 11:58 |
| | MB 240-207343/6 | 146535.D | 11/18/2015 | 12:20 |
| | LCS 240-207343/7 | 146536.D | 11/18/2015 | 12:41 |
| F-VS-UST-P-1 | 240-57930-1 | 146551.D | 11/18/2015 | 18:03 |
| F-VS-UST-P-3 | 240-57930-3 | 146552.D | 11/18/2015 | 18:25 |
| F-VB-UST-P-1 | 240-57930-5 | 146556.D | 11/18/2015 | 19:51 |
| F-VB-UST-P-2 | 240-57930-6 | 146557.D | 11/18/2015 | 20:12 |
| F-VB-UST-P-3 | 240-57930-7 | 146558.D | 11/18/2015 | 20:33 |
| F-VB-UST-P-4 | 240-57930-8 | 146559.D | 11/18/2015 | 20:55 |
| F-VB-UST-P-5 | 240-57930-9 | 146560.D | 11/18/2015 | 21:16 |
| F-VB-UST-P-6 | 240-57930-10 | 146561.D | 11/18/2015 | 21:38 |
| F-VS-UST-P-DUP | 240-57930-11 | 146562.D | 11/18/2015 | 21:59 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB14136.D  BFB Injection Date: 11/19/2015

Instrument ID: A3UX14  BFB Injection Time: 10:42

Analysis Batch No.: 207560

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 19.7 | |
| 75 | 30.0 - 60.0 % of mass 95 | 47.3 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.8 | |
| 173 | Less than 2.0 % of mass 174 | 0.8 | (1.0) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 78.8 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.9 | (7.4) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 75.4 | (95.6) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.6 | (6.1) 2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-207560/3 | 146565.D | 11/19/2015 | 11:25 |
| | CCV 240-207560/4 | 146566.D | 11/19/2015 | 11:46 |
| | LCS 240-207560/5 | 146567.D | 11/19/2015 | 12:08 |
| | MB 240-207560/6 | 146568.D | 11/19/2015 | 12:29 |
| | LCS 240-207560/7 | 146569.D | 11/19/2015 | 12:51 |
| F-VS-UST-P-4 | 240-57930-4 | 146581.D | 11/19/2015 | 17:08 |
| F-VS-UST-P-2 | 240-57930-2 | 146583.D | 11/19/2015 | 17:52 |
| F-VS-UST-P-2 | 240-57930-2 | 146584.D | 11/19/2015 | 18:13 |
| | 240-57921-A-4-J MS | 146590.D | 11/19/2015 | 20:22 |
| | 240-57921-A-4-K MSD | 146591.D | 11/19/2015 | 20:43 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB8851.D  BFB Injection Date: 07/15/2015

Instrument ID: A3UX8  BFB Injection Time: 00:55

Analysis Batch No.: 189147

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 32.0 | |
| 75 | 30.0 - 60.0 % of mass 95 | 57.2 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.5 | |
| 173 | Less than 2.0 % of mass 174 | 0.2 | (0.3) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 80.0 | |
| 175 | 5.0 - 9.0 % of mass 174 | 6.2 | (7.8) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 77.4 | (96.7) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 5.0 | (6.5) 2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STDA9 240-189147/4 | UX88545.D | 07/15/2015 | 01:51 |
| | STDA9 240-189147/5 | UX88546.D | 07/15/2015 | 02:12 |
| | STDA9 240-189147/6 | UX88547.D | 07/15/2015 | 02:34 |
| | STDA9 240-189147/7 | UX88548.D | 07/15/2015 | 02:55 |
| | STDA9 240-189147/8 | UX88549.D | 07/15/2015 | 03:16 |
| | STDA9 240-189147/9 | UX88550.D | 07/15/2015 | 03:38 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB8863.D  BFB Injection Date: 07/30/2015

Instrument ID: A3UX8  BFB Injection Time: 02:03

Analysis Batch No.: 191232

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 31.7 | |
| 75 | 30.0 - 60.0 % of mass 95 | 58.3 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.7 | |
| 173 | Less than 2.0 % of mass 174 | 0.0 | (0.0) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 79.4 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.7 | (7.2) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 77.2 | (97.3) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 5.1 | (6.6) 2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD8260 240-191232/3 | UX88718.D | 07/30/2015 | 02:35 |
| | STD8260 240-191232/4 | UX88719.D | 07/30/2015 | 02:56 |
| | STD8260 240-191232/5 | UX88720.D | 07/30/2015 | 03:17 |
| | ICIS 240-191232/6 | UX88721.D | 07/30/2015 | 03:39 |
| | STD8260 240-191232/7 | UX88722.D | 07/30/2015 | 04:00 |
| | STD8260 240-191232/8 | UX88723.D | 07/30/2015 | 04:21 |
| | STD8260 240-191232/9 | UX88724.D | 07/30/2015 | 04:42 |
| | STD8260 240-191232/10 | UX88725.D | 07/30/2015 | 05:04 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB8945.D BFB Injection Date: 11/17/2015

Instrument ID: A3UX8 BFB Injection Time: 03:23

Analysis Batch No.: 206828

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 26.8 | |
| 75 | 30.0 - 60.0 % of mass 95 | 51.0 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.4 | |
| 173 | Less than 2.0 % of mass 174 | 0.5 | (0.6) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 83.9 | |
| 175 | 5.0 - 9.0 % of mass 174 | 6.1 | (7.3) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 80.6 | (96.1) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 5.3 | (6.6) 2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD8260 240-206828/3 | UX80794.D | 11/17/2015 | 03:57 |
| | STD8260 240-206828/4 | UX80795.D | 11/17/2015 | 04:18 |
| | STD8260 240-206828/5 | UX80796.D | 11/17/2015 | 04:39 |
| | ICIS 240-206828/6 | UX80797.D | 11/17/2015 | 05:01 |
| | STD8260 240-206828/7 | UX80798.D | 11/17/2015 | 05:34 |
| | STD8260 240-206828/8 | UX80799.D | 11/17/2015 | 05:55 |
| | STD8260 240-206828/9 | UX80800.D | 11/17/2015 | 06:16 |
| | STD8260 240-206828/10 | UX80801.D | 11/17/2015 | 06:38 |
| | ICV 240-206828/11 | UX80802.D | 11/17/2015 | 06:59 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: BFB8948.D  BFB Injection Date: 11/20/2015

Instrument ID: A3UX8  BFB Injection Time: 03:12

Analysis Batch No.: 207720

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 26.3 | |
| 75 | 30.0 - 60.0 % of mass 95 | 49.9 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.8 | |
| 173 | Less than 2.0 % of mass 174 | 0.5 | (0.6) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 81.2 | |
| 175 | 5.0 - 9.0 % of mass 174 | 6.0 | (7.3) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 78.2 | (96.3) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 5.1 | (6.5) 2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-207720/4 | UX80875.D | 11/20/2015 | 03:43 |
| | CCV 240-207720/5 | UX80876.D | 11/20/2015 | 04:04 |
| | LCS 240-207720/6 | UX80877.D | 11/20/2015 | 04:26 |
| | MB 240-207720/8 | UX80879.D | 11/20/2015 | 05:08 |
| | 240-57972-C-3-C MS | UX80889.D | 11/20/2015 | 08:41 |
| | 240-57972-C-3-D MSD | UX80890.D | 11/20/2015 | 09:02 |
| F-VS-UST-P-2 | 240-57930-2 | UX80893.D | 11/20/2015 | 10:06 |

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: STD8260 240-206893/4  Date Analyzed: 11/16/2015 10:07

Instrument ID: A3UX10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UXX3041.D  Heated Purge: (Y/N) N

Calibration ID: 31566

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 1573848 | 5.54 | 1127244 | 8.23 | 662837 | 10.47 |
| UPPER LIMIT | | 3147696 | 6.04 | 2254488 | 8.73 | 1325674 | 10.97 |
| LOWER LIMIT | | 786924 | 5.04 | 563622 | 7.73 | 331419 | 9.97 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-206893/16 | | 1268597 | 5.54 | 936859 | 8.23 | 485454 | 10.47 |
| CCVIS 240-207096/2 | | 1260986 | 5.54 | 901961 | 8.23 | 487617 | 10.46 |
| MB 240-207096/6 | | 1120609 | 5.54 | 835430 | 8.23 | 430567 | 10.47 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCVIS 240-207096/2  Date Analyzed: 11/17/2015 10:50

Instrument ID: A3UX10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UXX3066.D  Heated Purge: (Y/N) N

Calibration ID: 31568

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1260986 | 5.54 | 901961 | 8.23 | 487617 | 10.46 |
| UPPER LIMIT | | 2521972 | 6.04 | 1803922 | 8.73 | 975234 | 10.96 |
| LOWER LIMIT | | 630493 | 5.04 | 450981 | 7.73 | 243809 | 9.96 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-207096/3 | | 1228989 | 5.54 | 899191 | 8.23 | 458011 | 10.47 |
| LCS 240-207096/4 | | 1225707 | 5.54 | 880516 | 8.23 | 490859 | 10.47 |
| MB 240-207096/6 | | 1120609 | 5.54 | 835430 | 8.23 | 430567 | 10.47 |
| 240-57930-12 | TB-111315 | 1004779 | 5.54 | 746987 | 8.23 | 396832 | 10.47 |
| 200-30727-B-1 MS | | 1136775 | 5.54 | 814318 | 8.23 | 471751 | 10.46 |
| 200-30727-B-1 MSD | | 1161812 | 5.54 | 855630 | 8.23 | 485911 | 10.46 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Sample No.: ICIS 240-196026/6 Date Analyzed: 09/02/2015 15:37

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm)

Lab File ID (Standard): 144938.D Heated Purge: (Y/N) Y

Calibration ID: 30511

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 1489699 | 6.64 | 910038 | 9.37 | 406298 | 11.34 |
| UPPER LIMIT | | 2979398 | 7.14 | 1820076 | 9.87 | 812596 | 11.84 |
| LOWER LIMIT | | 744850 | 6.14 | 455019 | 8.87 | 203149 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCVIS 240-207343/3 | | 1130206 | 6.64 | 702239 | 9.36 | 329914 | 11.34 |
| CCVIS 240-207560/3 | | 1296804 | 6.64 | 819718 | 9.37 | 379204 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCVIS 240-207343/3  Date Analyzed: 11/18/2015 11:15

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 146532.D  Heated Purge: (Y/N) Y

Calibration ID: 30511

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1130206 | 6.64 | 702239 | 9.36 | 329914 | 11.34 |
| UPPER LIMIT | | 2260412 | 7.14 | 1404478 | 9.86 | 659828 | 11.84 |
| LOWER LIMIT | | 565103 | 6.14 | 351120 | 8.86 | 164957 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-207343/4 | | 1217421 | 6.64 | 740840 | 9.36 | 344747 | 11.34 |
| LCS 240-207343/5 | | 1262829 | 6.64 | 755702 | 9.36 | 373208 | 11.34 |
| MB 240-207343/6 | | 1229280 | 6.64 | 750179 | 9.36 | 371381 | 11.34 |
| LCS 240-207343/7 | | 1257343 | 6.64 | 779199 | 9.36 | 370145 | 11.34 |
| 240-57930-1 | F-VS-UST-P-1 | 994924 | 6.64 | 574643 | 9.36 | 241066 | 11.34 |
| 240-57930-3 | F-VS-UST-P-3 | 1158152 | 6.64 | 664830 | 9.36 | 277450 | 11.34 |
| 240-57930-5 | F-VB-UST-P-1 | 1271214 | 6.64 | 763958 | 9.36 | 380397 | 11.34 |
| 240-57930-6 | F-VB-UST-P-2 | 1220498 | 6.64 | 748569 | 9.36 | 376645 | 11.34 |
| 240-57930-7 | F-VB-UST-P-3 | 1152061 | 6.64 | 716657 | 9.36 | 355890 | 11.34 |
| 240-57930-8 | F-VB-UST-P-4 | 1145728 | 6.64 | 720294 | 9.36 | 354172 | 11.34 |
| 240-57930-9 | F-VB-UST-P-5 | 1332976 | 6.64 | 804271 | 9.36 | 390098 | 11.34 |
| 240-57930-10 | F-VB-UST-P-6 | 1150668 | 6.64 | 708332 | 9.36 | 343986 | 11.34 |
| 240-57930-11 | F-VS-UST-P-DUP | 1144818 | 6.64 | 667602 | 9.36 | 303312 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCVIS 240-207560/3  Date Analyzed: 11/19/2015 11:25

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 146565.D  Heated Purge: (Y/N) Y

Calibration ID: 30511

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1296804 | 6.64 | 819718 | 9.37 | 379204 | 11.34 |
| UPPER LIMIT | | 2593608 | 7.14 | 1639436 | 9.87 | 758408 | 11.84 |
| LOWER LIMIT | | 648402 | 6.14 | 409859 | 8.87 | 189602 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-207560/4 | | 1173494 | 6.64 | 720210 | 9.36 | 344802 | 11.34 |
| LCS 240-207560/5 | | 1277930 | 6.64 | 777122 | 9.36 | 384697 | 11.34 |
| MB 240-207560/6 | | 1178385 | 6.64 | 719057 | 9.37 | 360387 | 11.34 |
| LCS 240-207560/7 | | 1286546 | 6.64 | 796373 | 9.36 | 388037 | 11.34 |
| 240-57930-4 | F-VS-UST-P-4 | 1153644 | 6.64 | 520223 | 9.36 | 164672* | 11.34 |
| 240-57930-2 | F-VS-UST-P-2 | 1286592 | 6.64 | 782678 | 9.36 | 367359 | 11.34 |
| 240-57930-2 | F-VS-UST-P-2 | 1196312 | 6.64 | 653872 | 9.36 | 237602 | 11.34 |
| 240-57921-A-4-J MS | | 1466811 | 6.64 | 901613 | 9.36 | 435827 | 11.34 |
| 240-57921-A-4-K MSD | | 1482454 | 6.64 | 917456 | 9.36 | 430677 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: ICIS 240-206828/6  Date Analyzed: 11/17/2015 05:01

Instrument ID: A3UX8  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UX80797.D  Heated Purge: (Y/N) Y

Calibration ID: 31578

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 1908665 | 5.16 | 1226226 | 7.69 | 459813 | 9.88 |
| UPPER LIMIT | | 3817330 | 5.66 | 2452452 | 8.19 | 919626 | 10.38 |
| LOWER LIMIT | | 954333 | 4.66 | 613113 | 7.19 | 229907 | 9.38 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-206828/11 | | 1732390 | 5.16 | 1085618 | 7.69 | 421779 | 9.88 |
| CCVIS 240-207720/4 | | 1766506 | 5.16 | 1084766 | 7.69 | 412435 | 9.88 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCVIS 240-207720/4  Date Analyzed: 11/20/2015 03:43

Instrument ID: A3UX8  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UX80875.D  Heated Purge: (Y/N) Y

Calibration ID: 31578

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1766506 | 5.16 | 1084766 | 7.69 | 412435 | 9.88 |
| UPPER LIMIT | | 3533012 | 5.66 | 2169532 | 8.19 | 824870 | 10.38 |
| LOWER LIMIT | | 883253 | 4.66 | 542383 | 7.19 | 206218 | 9.38 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-207720/5 | | 1999155 | 5.16 | 1398829 | 7.69 | 506592 | 9.88 |
| LCS 240-207720/6 | | 1683051 | 5.16 | 1054787 | 7.69 | 392198 | 9.88 |
| MB 240-207720/8 | | 1662330 | 5.16 | 1019809 | 7.69 | 403715 | 9.88 |
| 240-57972-C-3-C MS | | 1643283 | 5.16 | 1027699 | 7.69 | 399822 | 9.88 |
| 240-57972-C-3-D MSD | | 1870371 | 5.16 | 1265241 | 7.69 | 455695 | 9.88 |
| 240-57930-2 | F-VS-UST-P-2 | 1653633 | 5.16 | 983279 | 7.69 | 338150 | 9.88 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-1 | Lab Sample ID: 240-57930-1

Matrix: Solid | Lab File ID: 146551.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:30

Sample wt/vol: 5.18(g) | Date Analyzed: 11/18/2015 18:03

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.57 | U | 5.7 | 0.57 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.74 | U | 5.7 | 0.74 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.35 | U | 5.7 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.48 | U | 5.7 | 0.48 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.7 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.91 | U | 5.7 | 0.91 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 5.7 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.57 | U | 5.7 | 0.57 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.89 | U | 5.7 | 0.89 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.87 | J | 5.7 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.43 | U | 5.7 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.66 | U | 5.7 | 0.66 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 11 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.7 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.52 | U | 5.7 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.7 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.78 | U | 5.7 | 0.78 |
| 142-28-9 | 1,3-Dichloropropane | 0.45 | U | 5.7 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.7 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.50 | U | 5.7 | 0.50 |
| 78-93-3 | 2-Butanone (MEK) | 18 | J | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.73 | U | 57 | 0.73 |
| 95-49-8 | 2-Chlorotoluene | 0.34 | U | 5.7 | 0.34 |
| 591-78-6 | 2-Hexanone | 0.77 | U | 23 | 0.77 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.7 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.7 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 86 | | 23 | 6.5 |
| 71-43-2 | Benzene | 1.8 | J | 5.7 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.7 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.7 | 0.27 |
| 74-83-9 | Bromomethane | 0.48 | U | 5.7 | 0.48 |
| 75-15-0 | Carbon disulfide | 0.67 | U | 5.7 | 0.67 |
| 56-23-5 | Carbon tetrachloride | 0.74 | U | 5.7 | 0.74 |
| 108-90-7 | Chlorobenzene | 0.60 | U | 5.7 | 0.60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-1 | Lab Sample ID: 240-57930-1

Matrix: Solid | Lab File ID: 146551.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:30

Sample wt/vol: 5.18(g) | Date Analyzed: 11/18/2015 18:03

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.42 | U | 5.7 | 0.42 |
| 75-00-3 | Chloroethane | 0.50 | U | 5.7 | 0.50 |
| 67-66-3 | Chloroform | 0.41 | U | 5.7 | 0.41 |
| 74-87-3 | Chloromethane | 0.90 | U | 5.7 | 0.90 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.7 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.81 | U | 5.7 | 0.81 |
| 74-95-3 | Dibromomethane | 0.48 | U | 5.7 | 0.48 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.7 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.7 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.7 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.5 | U | 230 | 6.5 |
| 100-41-4 | Ethylbenzene | 0.36 | U | 5.7 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 0.45 | U | 5.7 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.67 | U | 11 | 0.67 |
| 98-82-8 | Isopropylbenzene | 14 | | 5.7 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.49 | U | 5.7 | 0.49 |
| 75-09-2 | Methylene Chloride | 0.84 | U | 5.7 | 0.84 |
| 179601-23-1 | m-Xylene & p-Xylene | 1.9 | J | 11 | 0.48 |
| 91-20-3 | Naphthalene | 1.1 | J | 5.7 | 0.40 |
| 104-51-8 | n-Butylbenzene | 7.1 | | 5.7 | 0.67 |
| 103-65-1 | N-Propylbenzene | 20 | | 5.7 | 0.68 |
| 95-47-6 | o-Xylene | 0.75 | J | 5.7 | 0.65 |
| 135-98-8 | sec-Butylbenzene | 8.6 | | 5.7 | 1.0 |
| 100-42-5 | Styrene | 0.45 | U | 5.7 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.49 | U | 5.7 | 0.49 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.7 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.91 | U | 5.7 | 0.91 |
| 108-88-3 | Toluene | 1.2 | J B | 5.7 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.36 | U | 5.7 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.7 | 0.33 |
| 79-01-6 | Trichloroethene | 0.43 | U | 5.7 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.7 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.75 | U | 11 | 0.75 |
| 75-01-4 | Vinyl chloride | 0.34 | U | 5.7 | 0.34 |
| 1330-20-7 | Xylenes, Total | 2.7 | J | 11 | 0.61 |
| 98-06-6 | tert-Butylbenzene | 2.3 | J | 5.7 | 0.76 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-1 | Lab Sample ID: 240-57930-1

Matrix: Solid | Lab File ID: 146551.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:30

Sample wt/vol: 5.18(g) | Date Analyzed: 11/18/2015 18:03

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.60 | U | 5.7 | 0.60 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 104 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 100 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 97 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 104 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: 146583.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:35

Sample wt/vol: 5.74(g) | Date Analyzed: 11/19/2015 17:52

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.52 | U | 5.2 | 0.52 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.67 | U | 5.2 | 0.67 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.32 | U | 5.2 | 0.32 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.43 | U | 5.2 | 0.43 |
| 75-34-3 | 1,1-Dichloroethane | 0.35 | U | 5.2 | 0.35 |
| 75-35-4 | 1,1-Dichloroethene | 0.82 | U | 5.2 | 0.82 |
| 563-58-6 | 1,1-Dichloropropene | 0.38 | U | 5.2 | 0.38 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.52 | U | 5.2 | 0.52 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.80 | U | 5.2 | 0.80 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 7.9 | | 5.2 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.39 | U | 5.2 | 0.39 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 12 | | 5.2 | 0.60 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.8 | U | 10 | 2.8 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.35 | U | 5.2 | 0.35 |
| 107-06-2 | 1,2-Dichloroethane | 0.47 | U | 5.2 | 0.47 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.2 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.71 | U | 5.2 | 0.71 |
| 142-28-9 | 1,3-Dichloropropane | 0.41 | U | 5.2 | 0.41 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.29 | U | 5.2 | 0.29 |
| 594-20-7 | 2,2-Dichloropropane | 0.45 | U | 5.2 | 0.45 |
| 78-93-3 | 2-Butanone (MEK) | 6.1 | J | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.66 | U | 52 | 0.66 |
| 95-49-8 | 2-Chlorotoluene | 0.31 | U | 5.2 | 0.31 |
| 591-78-6 | 2-Hexanone | 0.70 | U | 21 | 0.70 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.2 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.85 | J | 5.2 | 0.42 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 71-43-2 | Benzene | 0.93 | U | 5.2 | 0.93 |
| 108-86-1 | Bromobenzene | 0.34 | U | 5.2 | 0.34 |
| 75-25-2 | Bromoform | 0.25 | U | 5.2 | 0.25 |
| 74-83-9 | Bromomethane | 0.43 | U | 5.2 | 0.43 |
| 75-15-0 | Carbon disulfide | 0.61 | U | 5.2 | 0.61 |
| 56-23-5 | Carbon tetrachloride | 0.67 | U | 5.2 | 0.67 |
| 108-90-7 | Chlorobenzene | 0.55 | U | 5.2 | 0.55 |
| 124-48-1 | Dibromochloromethane | 0.38 | U | 5.2 | 0.38 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: 146583.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:35

Sample wt/vol: 5.74(g) | Date Analyzed: 11/19/2015 17:52

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 0.45 | U | 5.2 | 0.45 |
| 67-66-3 | Chloroform | 0.37 | U | 5.2 | 0.37 |
| 74-87-3 | Chloromethane | 0.81 | U | 5.2 | 0.81 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.29 | U | 5.2 | 0.29 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.73 | U | 5.2 | 0.73 |
| 74-95-3 | Dibromomethane | 0.43 | U | 5.2 | 0.43 |
| 74-97-5 | Bromochloromethane | 0.34 | U | 5.2 | 0.34 |
| 75-71-8 | Dichlorodifluoromethane | 0.29 | U | 5.2 | 0.29 |
| 75-27-4 | Bromodichloromethane | 0.30 | U | 5.2 | 0.30 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.9 | U | 210 | 5.9 |
| 100-41-4 | Ethylbenzene | 9.1 | | 5.2 | 0.33 |
| 87-68-3 | Hexachlorobutadiene | 0.41 | U | 5.2 | 0.41 |
| 108-20-3 | Diisopropyl ether | 0.61 | U | 10 | 0.61 |
| 98-82-8 | Isopropylbenzene | 8.5 | | 5.2 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.44 | U | 5.2 | 0.44 |
| 75-09-2 | Methylene Chloride | 0.76 | U | 5.2 | 0.76 |
| 179601-23-1 | m-Xylene & p-Xylene | 36 | | 10 | 0.43 |
| 91-20-3 | Naphthalene | 10 | | 5.2 | 0.36 |
| 104-51-8 | n-Butylbenzene | 4.6 | J | 5.2 | 0.61 |
| 103-65-1 | N-Propylbenzene | 9.7 | | 5.2 | 0.62 |
| 95-47-6 | o-Xylene | 2.7 | J | 5.2 | 0.59 |
| 135-98-8 | sec-Butylbenzene | 3.6 | J | 5.2 | 0.94 |
| 100-42-5 | Styrene | 0.41 | U | 5.2 | 0.41 |
| 994-05-8 | Tert-amyl methyl ether | 0.44 | U | 5.2 | 0.44 |
| 637-92-3 | Ethyl t-butyl ether | 0.42 | U | 5.2 | 0.42 |
| 127-18-4 | Tetrachloroethene | 0.82 | U | 5.2 | 0.82 |
| 108-88-3 | Toluene | 1.2 | J B | 5.2 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.33 | U | 5.2 | 0.33 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.30 | U | 5.2 | 0.30 |
| 79-01-6 | Trichloroethene | 0.39 | U | 5.2 | 0.39 |
| 75-69-4 | Trichlorofluoromethane | 0.30 | U | 5.2 | 0.30 |
| 108-05-4 | Vinyl acetate | 0.68 | U | 10 | 0.68 |
| 75-01-4 | Vinyl chloride | 0.31 | U | 5.2 | 0.31 |
| 1330-20-7 | Xylenes, Total | 39 | | 10 | 0.56 |
| 98-06-6 | tert-Butylbenzene | 0.87 | J | 5.2 | 0.69 |
| 106-93-4 | 1,2-Dibromoethane | 0.55 | U | 5.2 | 0.55 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 Lab Sample ID: 240-57930-2

Matrix: Solid Lab File ID: 146583.D

Analysis Method: 8260C Date Collected: 11/13/2015 10:35

Sample wt/vol: 5.74(g) Date Analyzed: 11/19/2015 17:52

Soil Aliquot Vol: Dilution Factor: 1

Soil Extract Vol.: GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 Level: (low/med) Low

Analysis Batch No.: 207560 Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 94 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 92 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 91 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 98 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: 146584.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:35

Sample wt/vol: 5.09(g) | Date Analyzed: 11/19/2015 18:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.76 | U | 5.8 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.91 | U | 5.8 | 0.91 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 98 | | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 170 | | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 5.8 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 8.8 | J | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 6.5 | | 5.8 | 0.48 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 71-43-2 | Benzene | 6.8 | | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.69 | U | 5.8 | 0.69 |
| 56-23-5 | Carbon tetrachloride | 0.76 | U | 5.8 | 0.76 |
| 108-90-7 | Chlorobenzene | 0.62 | U | 5.8 | 0.62 |
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: 146584.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:35

Sample wt/vol: 5.09(g) | Date Analyzed: 11/19/2015 18:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 5.8 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 5.8 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 140 | | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.69 | U | 12 | 0.69 |
| 98-82-8 | Isopropylbenzene | 63 | | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 520 | E | 12 | 0.49 |
| 91-20-3 | Naphthalene | 17 | | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 11 | | 5.8 | 0.69 |
| 103-65-1 | N-Propylbenzene | 55 | | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 14 | | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 12 | | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.48 | U | 5.8 | 0.48 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 4.0 | J B | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.77 | U | 12 | 0.77 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 530 | | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 4.4 | J | 5.8 | 0.78 |
| 106-93-4 | 1,2-Dibromoethane | 0.62 | U | 5.8 | 0.62 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: 146584.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:35

Sample wt/vol: 5.09(g) | Date Analyzed: 11/19/2015 18:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 103 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 92 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 105 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 95 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: UX80893.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:35

Sample wt/vol: 5.61(g) | Date Analyzed: 11/20/2015 10:06

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207720 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 67-64-1 | Acetone | 26 | B | 21 | 6.0 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 110 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 94 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 93 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 88 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-3 | Lab Sample ID: 240-57930-3

Matrix: Solid | Lab File ID: 146552.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:40

Sample wt/vol: 5.55(g) | Date Analyzed: 11/18/2015 18:25

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.53 | U | 5.3 | 0.53 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.69 | U | 5.3 | 0.69 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.33 | U | 5.3 | 0.33 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 5.3 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.36 | U | 5.3 | 0.36 |
| 75-35-4 | 1,1-Dichloroethene | 0.85 | U | 5.3 | 0.85 |
| 563-58-6 | 1,1-Dichloropropene | 0.39 | U | 5.3 | 0.39 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.53 | U | 5.3 | 0.53 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.83 | U | 5.3 | 0.83 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 1.5 | J | 5.3 | 0.49 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.40 | U | 5.3 | 0.40 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.6 | J | 5.3 | 0.62 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.9 | U | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.36 | U | 5.3 | 0.36 |
| 107-06-2 | 1,2-Dichloroethane | 0.49 | U | 5.3 | 0.49 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.3 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.73 | U | 5.3 | 0.73 |
| 142-28-9 | 1,3-Dichloropropane | 0.43 | U | 5.3 | 0.43 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.30 | U | 5.3 | 0.30 |
| 594-20-7 | 2,2-Dichloropropane | 0.47 | U | 5.3 | 0.47 |
| 78-93-3 | 2-Butanone (MEK) | 2.4 | J | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.68 | U | 53 | 0.68 |
| 95-49-8 | 2-Chlorotoluene | 0.32 | U | 5.3 | 0.32 |
| 591-78-6 | 2-Hexanone | 0.72 | U | 21 | 0.72 |
| 106-43-4 | 4-Chlorotoluene | 0.28 | U | 5.3 | 0.28 |
| 99-87-6 | p-Isopropyltoluene | 0.44 | U | 5.3 | 0.44 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 12 | J | 21 | 6.1 |
| 71-43-2 | Benzene | 0.96 | U | 5.3 | 0.96 |
| 108-86-1 | Bromobenzene | 0.35 | U | 5.3 | 0.35 |
| 75-25-2 | Bromoform | 0.26 | U | 5.3 | 0.26 |
| 74-83-9 | Bromomethane | 0.45 | U | 5.3 | 0.45 |
| 75-15-0 | Carbon disulfide | 0.63 | U | 5.3 | 0.63 |
| 56-23-5 | Carbon tetrachloride | 0.69 | U | 5.3 | 0.69 |
| 108-90-7 | Chlorobenzene | 0.56 | U | 5.3 | 0.56 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-3 | Lab Sample ID: 240-57930-3

Matrix: Solid | Lab File ID: 146552.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:40

Sample wt/vol: 5.55(g) | Date Analyzed: 11/18/2015 18:25

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.39 | U | 5.3 | 0.39 |
| 75-00-3 | Chloroethane | 0.47 | U | 5.3 | 0.47 |
| 67-66-3 | Chloroform | 0.38 | U | 5.3 | 0.38 |
| 74-87-3 | Chloromethane | 0.84 | U | 5.3 | 0.84 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.30 | U | 5.3 | 0.30 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.76 | U | 5.3 | 0.76 |
| 74-95-3 | Dibromomethane | 0.45 | U | 5.3 | 0.45 |
| 74-97-5 | Bromochloromethane | 0.35 | U | 5.3 | 0.35 |
| 75-71-8 | Dichlorodifluoromethane | 0.30 | U | 5.3 | 0.30 |
| 75-27-4 | Bromodichloromethane | 0.31 | U | 5.3 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.1 | U | 210 | 6.1 |
| 100-41-4 | Ethylbenzene | 0.34 | U | 5.3 | 0.34 |
| 87-68-3 | Hexachlorobutadiene | 0.43 | U | 5.3 | 0.43 |
| 108-20-3 | Diisopropyl ether | 0.63 | U | 11 | 0.63 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.3 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.46 | U | 5.3 | 0.46 |
| 75-09-2 | Methylene Chloride | 0.79 | U | 5.3 | 0.79 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.45 | U | 11 | 0.45 |
| 91-20-3 | Naphthalene | 1.7 | J | 5.3 | 0.37 |
| 104-51-8 | n-Butylbenzene | 0.63 | U | 5.3 | 0.63 |
| 103-65-1 | N-Propylbenzene | 0.64 | U | 5.3 | 0.64 |
| 95-47-6 | o-Xylene | 0.61 | U | 5.3 | 0.61 |
| 135-98-8 | sec-Butylbenzene | 0.97 | U | 5.3 | 0.97 |
| 100-42-5 | Styrene | 0.43 | U | 5.3 | 0.43 |
| 994-05-8 | Tert-amyl methyl ether | 0.46 | U | 5.3 | 0.46 |
| 637-92-3 | Ethyl t-butyl ether | 0.44 | U | 5.3 | 0.44 |
| 127-18-4 | Tetrachloroethene | 0.85 | U | 5.3 | 0.85 |
| 108-88-3 | Toluene | 0.43 | J B | 5.3 | 0.29 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.34 | U | 5.3 | 0.34 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.31 | U | 5.3 | 0.31 |
| 79-01-6 | Trichloroethene | 0.40 | U | 5.3 | 0.40 |
| 75-69-4 | Trichlorofluoromethane | 0.31 | U | 5.3 | 0.31 |
| 108-05-4 | Vinyl acetate | 0.70 | U | 11 | 0.70 |
| 75-01-4 | Vinyl chloride | 0.32 | U | 5.3 | 0.32 |
| 1330-20-7 | Xylenes, Total | 0.57 | U | 11 | 0.57 |
| 98-06-6 | tert-Butylbenzene | 0.71 | U | 5.3 | 0.71 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-3 | Lab Sample ID: 240-57930-3

Matrix: Solid | Lab File ID: 146552.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:40

Sample wt/vol: 5.55(g) | Date Analyzed: 11/18/2015 18:25

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.56 | U | 5.3 | 0.56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 94 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 91 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 89 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 91 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-4 | Lab Sample ID: 240-57930-4

Matrix: Solid | Lab File ID: 146581.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:45

Sample wt/vol: 5.24(g) | Date Analyzed: 11/19/2015 17:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.9 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.57 | U | 5.7 | 0.57 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.7 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U * | 5.7 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.48 | U | 5.7 | 0.48 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.7 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.92 | U | 5.7 | 0.92 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 5.7 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.57 | U * | 5.7 | 0.57 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U * | 5.7 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U * | 5.7 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U * | 5.7 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U * | 5.7 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U * | 11 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U * | 5.7 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.7 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.7 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.79 | U * | 5.7 | 0.79 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.7 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U * | 5.7 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.7 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.4 | J | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.73 | U | 57 | 0.73 |
| 95-49-8 | 2-Chlorotoluene | 0.34 | U * | 5.7 | 0.34 |
| 591-78-6 | 2-Hexanone | 0.78 | U | 23 | 0.78 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U * | 5.7 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U * | 5.7 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 21 | J | 23 | 6.5 |
| 71-43-2 | Benzene | 1.0 | U | 5.7 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U * | 5.7 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.7 | 0.28 |
| 74-83-9 | Bromomethane | 0.48 | U | 5.7 | 0.48 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.7 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.7 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.7 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-4 | Lab Sample ID: 240-57930-4

Matrix: Solid | Lab File ID: 146581.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:45

Sample wt/vol: 5.24(g) | Date Analyzed: 11/19/2015 17:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.9 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.42 | U | 5.7 | 0.42 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.7 | 0.51 |
| 67-66-3 | Chloroform | 0.41 | U | 5.7 | 0.41 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.7 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.7 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.7 | 0.82 |
| 74-95-3 | Dibromomethane | 0.48 | U | 5.7 | 0.48 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.7 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.7 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.7 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.7 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U * | 5.7 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 11 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.7 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.49 | U | 5.7 | 0.49 |
| 75-09-2 | Methylene Chloride | 0.85 | U | 5.7 | 0.85 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.48 | U | 11 | 0.48 |
| 91-20-3 | Naphthalene | 0.88 | J * | 5.7 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U * | 5.7 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U * | 5.7 | 0.69 |
| 95-47-6 | o-Xylene | 0.65 | U | 5.7 | 0.65 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U * | 5.7 | 1.0 |
| 100-42-5 | Styrene | 0.46 | U | 5.7 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.49 | U | 5.7 | 0.49 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.7 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.92 | U | 5.7 | 0.92 |
| 108-88-3 | Toluene | 0.47 | J B | 5.7 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.7 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.7 | 0.33 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.7 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.7 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 11 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.34 | U | 5.7 | 0.34 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 11 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U * | 5.7 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-4 | Lab Sample ID: 240-57930-4

Matrix: Solid | Lab File ID: 146581.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:45

Sample wt/vol: 5.24(g) | Date Analyzed: 11/19/2015 17:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.9 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.7 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 111 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 87 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 102 | * | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 85 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 | Lab Sample ID: 240-57930-5

Matrix: Solid | Lab File ID: 146556.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:00

Sample wt/vol: 5.06(g) | Date Analyzed: 11/18/2015 19:51

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.76 | U | 5.8 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 5.8 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.91 | U | 5.8 | 0.91 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.54 | U | 5.8 | 0.54 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.68 | U | 5.8 | 0.68 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 5.8 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.54 | U | 5.8 | 0.54 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.81 | U | 5.8 | 0.81 |
| 142-28-9 | 1,3-Dichloropropane | 0.47 | U | 5.8 | 0.47 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 5.8 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 1.3 | J | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.75 | U | 58 | 0.75 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.48 | U | 5.8 | 0.48 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 8.2 | J | 23 | 6.7 |
| 71-43-2 | Benzene | 1.1 | U | 5.8 | 1.1 |
| 108-86-1 | Bromobenzene | 0.39 | U | 5.8 | 0.39 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.69 | U | 5.8 | 0.69 |
| 56-23-5 | Carbon tetrachloride | 0.76 | U | 5.8 | 0.76 |
| 108-90-7 | Chlorobenzene | 0.62 | U | 5.8 | 0.62 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 | Lab Sample ID: 240-57930-5

Matrix: Solid | Lab File ID: 146556.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:00

Sample wt/vol: 5.06(g) | Date Analyzed: 11/18/2015 19:51

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 5.8 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.83 | U | 5.8 | 0.83 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.39 | U | 5.8 | 0.39 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 5.8 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.7 | U | 230 | 6.7 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.47 | U | 5.8 | 0.47 |
| 108-20-3 | Diisopropyl ether | 0.69 | U | 12 | 0.69 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.41 | U | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 0.69 | U | 5.8 | 0.69 |
| 103-65-1 | N-Propylbenzene | 0.70 | U | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 0.67 | U | 5.8 | 0.67 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.47 | U | 5.8 | 0.47 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.48 | U | 5.8 | 0.48 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.32 | U | 5.8 | 0.32 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.77 | U | 12 | 0.77 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.63 | U | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 0.78 | U | 5.8 | 0.78 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 | Lab Sample ID: 240-57930-5

Matrix: Solid | Lab File ID: 146556.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:00

Sample wt/vol: 5.06(g) | Date Analyzed: 11/18/2015 19:51

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.62 | U | 5.8 | 0.62 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 89 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 78 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 90 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-2 | Lab Sample ID: 240-57930-6

Matrix: Solid | Lab File ID: 146557.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:05

Sample wt/vol: 5.31(g) | Date Analyzed: 11/18/2015 20:12

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.5 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.56 | U | 5.6 | 0.56 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.72 | U | 5.6 | 0.72 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.35 | U | 5.6 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.47 | U | 5.6 | 0.47 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.6 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.89 | U | 5.6 | 0.89 |
| 563-58-6 | 1,1-Dichloropropene | 0.41 | U | 5.6 | 0.41 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.56 | U | 5.6 | 0.56 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.87 | U | 5.6 | 0.87 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.51 | U | 5.6 | 0.51 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.42 | U | 5.6 | 0.42 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.65 | U | 5.6 | 0.65 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.6 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.51 | U | 5.6 | 0.51 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.6 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.77 | U | 5.6 | 0.77 |
| 142-28-9 | 1,3-Dichloropropane | 0.45 | U | 5.6 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.6 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.6 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 5.4 | J | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.71 | U | 56 | 0.71 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.6 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.76 | U | 22 | 0.76 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.6 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.46 | U | 5.6 | 0.46 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 70 | | 22 | 6.4 |
| 71-43-2 | Benzene | 1.0 | U | 5.6 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.6 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.6 | 0.27 |
| 74-83-9 | Bromomethane | 0.47 | U | 5.6 | 0.47 |
| 75-15-0 | Carbon disulfide | 0.66 | U | 5.6 | 0.66 |
| 56-23-5 | Carbon tetrachloride | 0.72 | U | 5.6 | 0.72 |
| 108-90-7 | Chlorobenzene | 0.59 | U | 5.6 | 0.59 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-2 | Lab Sample ID: 240-57930-6

Matrix: Solid | Lab File ID: 146557.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:05

Sample wt/vol: 5.31(g) | Date Analyzed: 11/18/2015 20:12

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.5 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.41 | U | 5.6 | 0.41 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.6 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.6 | 0.40 |
| 74-87-3 | Chloromethane | 0.88 | U | 5.6 | 0.88 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.6 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.79 | U | 5.6 | 0.79 |
| 74-95-3 | Dibromomethane | 0.47 | U | 5.6 | 0.47 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.6 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.6 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.6 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.4 | U | 220 | 6.4 |
| 100-41-4 | Ethylbenzene | 0.36 | U | 5.6 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 0.45 | U | 5.6 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.66 | U | 11 | 0.66 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.6 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.6 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.82 | U | 5.6 | 0.82 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.47 | U | 11 | 0.47 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.6 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.66 | U | 5.6 | 0.66 |
| 103-65-1 | N-Propylbenzene | 0.67 | U | 5.6 | 0.67 |
| 95-47-6 | o-Xylene | 0.64 | U | 5.6 | 0.64 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.6 | 1.0 |
| 100-42-5 | Styrene | 0.45 | U | 5.6 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.6 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.46 | U | 5.6 | 0.46 |
| 127-18-4 | Tetrachloroethene | 0.89 | U | 5.6 | 0.89 |
| 108-88-3 | Toluene | 0.30 | U | 5.6 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.36 | U | 5.6 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.6 | 0.32 |
| 79-01-6 | Trichloroethene | 0.42 | U | 5.6 | 0.42 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.6 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.74 | U | 11 | 0.74 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.6 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.60 | U | 11 | 0.60 |
| 98-06-6 | tert-Butylbenzene | 0.75 | U | 5.6 | 0.75 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-2 | Lab Sample ID: 240-57930-6

Matrix: Solid | Lab File ID: 146557.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:05

Sample wt/vol: 5.31(g) | Date Analyzed: 11/18/2015 20:12

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.5 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.59 | U | 5.6 | 0.59 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 95 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 79 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 100 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-3 | Lab Sample ID: 240-57930-7

Matrix: Solid | Lab File ID: 146558.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:10

Sample wt/vol: 5.29(g) | Date Analyzed: 11/18/2015 20:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.0 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.76 | U | 5.8 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 5.8 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.91 | U | 5.8 | 0.91 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.54 | U | 5.8 | 0.54 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.68 | U | 5.8 | 0.68 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 5.8 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.54 | U | 5.8 | 0.54 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.81 | U | 5.8 | 0.81 |
| 142-28-9 | 1,3-Dichloropropane | 0.47 | U | 5.8 | 0.47 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 5.8 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.9 | J | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.75 | U | 58 | 0.75 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.48 | U | 5.8 | 0.48 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 66 | | 23 | 6.7 |
| 71-43-2 | Benzene | 1.1 | U | 5.8 | 1.1 |
| 108-86-1 | Bromobenzene | 0.39 | U | 5.8 | 0.39 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.69 | U | 5.8 | 0.69 |
| 56-23-5 | Carbon tetrachloride | 0.76 | U | 5.8 | 0.76 |
| 108-90-7 | Chlorobenzene | 0.62 | U | 5.8 | 0.62 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-3 | Lab Sample ID: 240-57930-7

Matrix: Solid | Lab File ID: 146558.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:10

Sample wt/vol: 5.29(g) | Date Analyzed: 11/18/2015 20:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.0 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 5.8 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.83 | U | 5.8 | 0.83 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.39 | U | 5.8 | 0.39 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 5.8 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.7 | U | 230 | 6.7 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.47 | U | 5.8 | 0.47 |
| 108-20-3 | Diisopropyl ether | 0.69 | U | 12 | 0.69 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.41 | U | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 0.69 | U | 5.8 | 0.69 |
| 103-65-1 | N-Propylbenzene | 0.70 | U | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 0.67 | U | 5.8 | 0.67 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.47 | U | 5.8 | 0.47 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.48 | U | 5.8 | 0.48 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.32 | U | 5.8 | 0.32 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.77 | U | 12 | 0.77 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.63 | U | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 0.78 | U | 5.8 | 0.78 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-3 | Lab Sample ID: 240-57930-7

Matrix: Solid | Lab File ID: 146558.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:10

Sample wt/vol: 5.29(g) | Date Analyzed: 11/18/2015 20:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.0 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.62 | U | 5.8 | 0.62 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 94 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 99 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 82 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 105 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-4 | Lab Sample ID: 240-57930-8

Matrix: Solid | Lab File ID: 146559.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:15

Sample wt/vol: 5.55(g) | Date Analyzed: 11/18/2015 20:55

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.1 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.54 | U | 5.4 | 0.54 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.70 | U | 5.4 | 0.70 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.33 | U | 5.4 | 0.33 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 5.4 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.37 | U | 5.4 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 0.86 | U | 5.4 | 0.86 |
| 563-58-6 | 1,1-Dichloropropene | 0.40 | U | 5.4 | 0.40 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.54 | U | 5.4 | 0.54 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.84 | U | 5.4 | 0.84 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.49 | U | 5.4 | 0.49 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.41 | U | 5.4 | 0.41 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.62 | U | 5.4 | 0.62 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.9 | U | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.37 | U | 5.4 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 0.49 | U | 5.4 | 0.49 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.4 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.74 | U | 5.4 | 0.74 |
| 142-28-9 | 1,3-Dichloropropane | 0.43 | U | 5.4 | 0.43 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.30 | U | 5.4 | 0.30 |
| 594-20-7 | 2,2-Dichloropropane | 0.47 | U | 5.4 | 0.47 |
| 78-93-3 | 2-Butanone (MEK) | 1.9 | J | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.69 | U | 54 | 0.69 |
| 95-49-8 | 2-Chlorotoluene | 0.32 | U | 5.4 | 0.32 |
| 591-78-6 | 2-Hexanone | 0.73 | U | 21 | 0.73 |
| 106-43-4 | 4-Chlorotoluene | 0.28 | U | 5.4 | 0.28 |
| 99-87-6 | p-Isopropyltoluene | 0.44 | U | 5.4 | 0.44 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 16 | J | 21 | 6.1 |
| 71-43-2 | Benzene | 0.97 | U | 5.4 | 0.97 |
| 108-86-1 | Bromobenzene | 0.35 | U | 5.4 | 0.35 |
| 75-25-2 | Bromoform | 0.26 | U | 5.4 | 0.26 |
| 74-83-9 | Bromomethane | 0.45 | U | 5.4 | 0.45 |
| 75-15-0 | Carbon disulfide | 0.63 | U | 5.4 | 0.63 |
| 56-23-5 | Carbon tetrachloride | 0.70 | U | 5.4 | 0.70 |
| 108-90-7 | Chlorobenzene | 0.57 | U | 5.4 | 0.57 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-4 | Lab Sample ID: 240-57930-8

Matrix: Solid | Lab File ID: 146559.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:15

Sample wt/vol: 5.55(g) | Date Analyzed: 11/18/2015 20:55

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.1 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.40 | U | 5.4 | 0.40 |
| 75-00-3 | Chloroethane | 0.47 | U | 5.4 | 0.47 |
| 67-66-3 | Chloroform | 0.39 | U | 5.4 | 0.39 |
| 74-87-3 | Chloromethane | 0.85 | U | 5.4 | 0.85 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.30 | U | 5.4 | 0.30 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.76 | U | 5.4 | 0.76 |
| 74-95-3 | Dibromomethane | 0.45 | U | 5.4 | 0.45 |
| 74-97-5 | Bromochloromethane | 0.35 | U | 5.4 | 0.35 |
| 75-71-8 | Dichlorodifluoromethane | 0.30 | U | 5.4 | 0.30 |
| 75-27-4 | Bromodichloromethane | 0.31 | U | 5.4 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.1 | U | 210 | 6.1 |
| 100-41-4 | Ethylbenzene | 0.34 | U | 5.4 | 0.34 |
| 87-68-3 | Hexachlorobutadiene | 0.43 | U | 5.4 | 0.43 |
| 108-20-3 | Diisopropyl ether | 0.63 | U | 11 | 0.63 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.4 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.46 | U | 5.4 | 0.46 |
| 75-09-2 | Methylene Chloride | 0.79 | U | 5.4 | 0.79 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.45 | U | 11 | 0.45 |
| 91-20-3 | Naphthalene | 0.38 | U | 5.4 | 0.38 |
| 104-51-8 | n-Butylbenzene | 0.63 | U | 5.4 | 0.63 |
| 103-65-1 | N-Propylbenzene | 0.64 | U | 5.4 | 0.64 |
| 95-47-6 | o-Xylene | 0.61 | U | 5.4 | 0.61 |
| 135-98-8 | sec-Butylbenzene | 0.98 | U | 5.4 | 0.98 |
| 100-42-5 | Styrene | 0.43 | U | 5.4 | 0.43 |
| 994-05-8 | Tert-amyl methyl ether | 0.46 | U | 5.4 | 0.46 |
| 637-92-3 | Ethyl t-butyl ether | 0.44 | U | 5.4 | 0.44 |
| 127-18-4 | Tetrachloroethene | 0.86 | U | 5.4 | 0.86 |
| 108-88-3 | Toluene | 0.29 | U | 5.4 | 0.29 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.34 | U | 5.4 | 0.34 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.31 | U | 5.4 | 0.31 |
| 79-01-6 | Trichloroethene | 0.41 | U | 5.4 | 0.41 |
| 75-69-4 | Trichlorofluoromethane | 0.31 | U | 5.4 | 0.31 |
| 108-05-4 | Vinyl acetate | 0.71 | U | 11 | 0.71 |
| 75-01-4 | Vinyl chloride | 0.32 | U | 5.4 | 0.32 |
| 1330-20-7 | Xylenes, Total | 0.58 | U | 11 | 0.58 |
| 98-06-6 | tert-Butylbenzene | 0.72 | U | 5.4 | 0.72 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-4 | Lab Sample ID: 240-57930-8

Matrix: Solid | Lab File ID: 146559.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:15

Sample wt/vol: 5.55(g) | Date Analyzed: 11/18/2015 20:55

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.1 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.57 | U | 5.4 | 0.57 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 96 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 103 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 89 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 113 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-5 | Lab Sample ID: 240-57930-9

Matrix: Solid | Lab File ID: 146560.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:20

Sample wt/vol: 5.55(g) | Date Analyzed: 11/18/2015 21:16

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.7 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.54 | U | 5.4 | 0.54 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.70 | U | 5.4 | 0.70 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.4 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 5.4 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.37 | U | 5.4 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 0.87 | U | 5.4 | 0.87 |
| 563-58-6 | 1,1-Dichloropropene | 0.40 | U | 5.4 | 0.40 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.54 | U | 5.4 | 0.54 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.84 | U | 5.4 | 0.84 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.50 | U | 5.4 | 0.50 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.41 | U | 5.4 | 0.41 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.63 | U | 5.4 | 0.63 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.9 | U | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.37 | U | 5.4 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 0.50 | U | 5.4 | 0.50 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.4 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 5.4 | 0.75 |
| 142-28-9 | 1,3-Dichloropropane | 0.43 | U | 5.4 | 0.43 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.30 | U | 5.4 | 0.30 |
| 594-20-7 | 2,2-Dichloropropane | 0.48 | U | 5.4 | 0.48 |
| 78-93-3 | 2-Butanone (MEK) | 2.0 | J | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.69 | U | 54 | 0.69 |
| 95-49-8 | 2-Chlorotoluene | 0.32 | U | 5.4 | 0.32 |
| 591-78-6 | 2-Hexanone | 0.74 | U | 22 | 0.74 |
| 106-43-4 | 4-Chlorotoluene | 0.28 | U | 5.4 | 0.28 |
| 99-87-6 | p-Isopropyltoluene | 0.44 | U | 5.4 | 0.44 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 22 | 1.3 |
| 67-64-1 | Acetone | 11 | J | 22 | 6.2 |
| 71-43-2 | Benzene | 0.97 | U | 5.4 | 0.97 |
| 108-86-1 | Bromobenzene | 0.36 | U | 5.4 | 0.36 |
| 75-25-2 | Bromoform | 0.26 | U | 5.4 | 0.26 |
| 74-83-9 | Bromomethane | 0.45 | U | 5.4 | 0.45 |
| 75-15-0 | Carbon disulfide | 0.64 | U | 5.4 | 0.64 |
| 56-23-5 | Carbon tetrachloride | 0.70 | U | 5.4 | 0.70 |
| 108-90-7 | Chlorobenzene | 0.57 | U | 5.4 | 0.57 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-5 | Lab Sample ID: 240-57930-9

Matrix: Solid | Lab File ID: 146560.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:20

Sample wt/vol: 5.55(g) | Date Analyzed: 11/18/2015 21:16

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.7 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.40 | U | 5.4 | 0.40 |
| 75-00-3 | Chloroethane | 0.48 | U | 5.4 | 0.48 |
| 67-66-3 | Chloroform | 0.39 | U | 5.4 | 0.39 |
| 74-87-3 | Chloromethane | 0.85 | U | 5.4 | 0.85 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.30 | U | 5.4 | 0.30 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.77 | U | 5.4 | 0.77 |
| 74-95-3 | Dibromomethane | 0.45 | U | 5.4 | 0.45 |
| 74-97-5 | Bromochloromethane | 0.36 | U | 5.4 | 0.36 |
| 75-71-8 | Dichlorodifluoromethane | 0.30 | U | 5.4 | 0.30 |
| 75-27-4 | Bromodichloromethane | 0.31 | U | 5.4 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.2 | U | 220 | 6.2 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.4 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.43 | U | 5.4 | 0.43 |
| 108-20-3 | Diisopropyl ether | 0.64 | U | 11 | 0.64 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.4 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.47 | U | 5.4 | 0.47 |
| 75-09-2 | Methylene Chloride | 0.80 | U | 5.4 | 0.80 |
| 179601-23-1 | m-Xylene & p-Xylene | 1.4 | J | 11 | 0.45 |
| 91-20-3 | Naphthalene | 0.38 | U | 5.4 | 0.38 |
| 104-51-8 | n-Butylbenzene | 0.64 | U | 5.4 | 0.64 |
| 103-65-1 | N-Propylbenzene | 0.65 | U | 5.4 | 0.65 |
| 95-47-6 | o-Xylene | 0.62 | U | 5.4 | 0.62 |
| 135-98-8 | sec-Butylbenzene | 0.98 | U | 5.4 | 0.98 |
| 100-42-5 | Styrene | 0.43 | U | 5.4 | 0.43 |
| 994-05-8 | Tert-amyl methyl ether | 0.47 | U | 5.4 | 0.47 |
| 637-92-3 | Ethyl t-butyl ether | 0.44 | U | 5.4 | 0.44 |
| 127-18-4 | Tetrachloroethene | 0.87 | U | 5.4 | 0.87 |
| 108-88-3 | Toluene | 0.75 | J B | 5.4 | 0.29 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.4 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.31 | U | 5.4 | 0.31 |
| 79-01-6 | Trichloroethene | 0.41 | U | 5.4 | 0.41 |
| 75-69-4 | Trichlorofluoromethane | 0.31 | U | 5.4 | 0.31 |
| 108-05-4 | Vinyl acetate | 0.71 | U | 11 | 0.71 |
| 75-01-4 | Vinyl chloride | 0.32 | U | 5.4 | 0.32 |
| 1330-20-7 | Xylenes, Total | 1.4 | J | 11 | 0.58 |
| 98-06-6 | tert-Butylbenzene | 0.72 | U | 5.4 | 0.72 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-5 | Lab Sample ID: 240-57930-9

Matrix: Solid | Lab File ID: 146560.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:20

Sample wt/vol: 5.55(g) | Date Analyzed: 11/18/2015 21:16

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.7 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.57 | U | 5.4 | 0.57 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 84 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 86 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 76 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 89 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-6 | Lab Sample ID: 240-57930-10

Matrix: Solid | Lab File ID: 146561.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:25

Sample wt/vol: 5.10(g) | Date Analyzed: 11/18/2015 21:38

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.1 | J | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 57 | | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-6 | Lab Sample ID: 240-57930-10

Matrix: Solid | Lab File ID: 146561.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:25

Sample wt/vol: 5.10(g) | Date Analyzed: 11/18/2015 21:38

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.40 | U | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-6 | Lab Sample ID: 240-57930-10

Matrix: Solid | Lab File ID: 146561.D

Analysis Method: 8260C | Date Collected: 11/13/2015 10:25

Sample wt/vol: 5.10(g) | Date Analyzed: 11/18/2015 21:38

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 93 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 97 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 86 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 105 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP | Lab Sample ID: 240-57930-11

Matrix: Solid | Lab File ID: 146562.D

Analysis Method: 8260C | Date Collected: 11/13/2015 00:00

Sample wt/vol: 5.73(g) | Date Analyzed: 11/18/2015 21:59

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.52 | U | 5.2 | 0.52 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.67 | U | 5.2 | 0.67 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.32 | U | 5.2 | 0.32 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.43 | U | 5.2 | 0.43 |
| 75-34-3 | 1,1-Dichloroethane | 0.35 | U | 5.2 | 0.35 |
| 75-35-4 | 1,1-Dichloroethene | 0.83 | U | 5.2 | 0.83 |
| 563-58-6 | 1,1-Dichloropropene | 0.38 | U | 5.2 | 0.38 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.52 | U | 5.2 | 0.52 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.81 | U | 5.2 | 0.81 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.2 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.39 | U | 5.2 | 0.39 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.60 | U | 5.2 | 0.60 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.8 | U | 10 | 2.8 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.35 | U | 5.2 | 0.35 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.2 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.2 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.71 | U | 5.2 | 0.71 |
| 142-28-9 | 1,3-Dichloropropane | 0.41 | U | 5.2 | 0.41 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.29 | U | 5.2 | 0.29 |
| 594-20-7 | 2,2-Dichloropropane | 0.45 | U | 5.2 | 0.45 |
| 78-93-3 | 2-Butanone (MEK) | 6.4 | J | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.66 | U | 52 | 0.66 |
| 95-49-8 | 2-Chlorotoluene | 0.31 | U | 5.2 | 0.31 |
| 591-78-6 | 2-Hexanone | 0.70 | U | 21 | 0.70 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.2 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.42 | U | 5.2 | 0.42 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 34 | | 21 | 5.9 |
| 71-43-2 | Benzene | 0.93 | U | 5.2 | 0.93 |
| 108-86-1 | Bromobenzene | 0.34 | U | 5.2 | 0.34 |
| 75-25-2 | Bromoform | 0.25 | U | 5.2 | 0.25 |
| 74-83-9 | Bromomethane | 0.43 | U | 5.2 | 0.43 |
| 75-15-0 | Carbon disulfide | 0.61 | U | 5.2 | 0.61 |
| 56-23-5 | Carbon tetrachloride | 0.67 | U | 5.2 | 0.67 |
| 108-90-7 | Chlorobenzene | 0.55 | U | 5.2 | 0.55 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP | Lab Sample ID: 240-57930-11

Matrix: Solid | Lab File ID: 146562.D

Analysis Method: 8260C | Date Collected: 11/13/2015 00:00

Sample wt/vol: 5.73(g) | Date Analyzed: 11/18/2015 21:59

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.38 | U | 5.2 | 0.38 |
| 75-00-3 | Chloroethane | 0.45 | U | 5.2 | 0.45 |
| 67-66-3 | Chloroform | 0.37 | U | 5.2 | 0.37 |
| 74-87-3 | Chloromethane | 0.82 | U | 5.2 | 0.82 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.29 | U | 5.2 | 0.29 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.73 | U | 5.2 | 0.73 |
| 74-95-3 | Dibromomethane | 0.43 | U | 5.2 | 0.43 |
| 74-97-5 | Bromochloromethane | 0.34 | U | 5.2 | 0.34 |
| 75-71-8 | Dichlorodifluoromethane | 0.29 | U | 5.2 | 0.29 |
| 75-27-4 | Bromodichloromethane | 0.30 | U | 5.2 | 0.30 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.9 | U | 210 | 5.9 |
| 100-41-4 | Ethylbenzene | 0.33 | U | 5.2 | 0.33 |
| 87-68-3 | Hexachlorobutadiene | 0.41 | U | 5.2 | 0.41 |
| 108-20-3 | Diisopropyl ether | 0.61 | U | 10 | 0.61 |
| 98-82-8 | Isopropylbenzene | 2.8 | J | 5.2 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.44 | U | 5.2 | 0.44 |
| 75-09-2 | Methylene Chloride | 0.76 | U | 5.2 | 0.76 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.43 | U | 10 | 0.43 |
| 91-20-3 | Naphthalene | 0.36 | U | 5.2 | 0.36 |
| 104-51-8 | n-Butylbenzene | 1.7 | J | 5.2 | 0.61 |
| 103-65-1 | N-Propylbenzene | 4.4 | J | 5.2 | 0.62 |
| 95-47-6 | o-Xylene | 0.59 | U | 5.2 | 0.59 |
| 135-98-8 | sec-Butylbenzene | 2.0 | J | 5.2 | 0.94 |
| 100-42-5 | Styrene | 0.41 | U | 5.2 | 0.41 |
| 994-05-8 | Tert-amyl methyl ether | 0.44 | U | 5.2 | 0.44 |
| 637-92-3 | Ethyl t-butyl ether | 0.42 | U | 5.2 | 0.42 |
| 127-18-4 | Tetrachloroethene | 0.83 | U | 5.2 | 0.83 |
| 108-88-3 | Toluene | 0.31 | J B | 5.2 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.33 | U | 5.2 | 0.33 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.30 | U | 5.2 | 0.30 |
| 79-01-6 | Trichloroethene | 0.39 | U | 5.2 | 0.39 |
| 75-69-4 | Trichlorofluoromethane | 0.30 | U | 5.2 | 0.30 |
| 108-05-4 | Vinyl acetate | 0.68 | U | 10 | 0.68 |
| 75-01-4 | Vinyl chloride | 0.31 | U | 5.2 | 0.31 |
| 1330-20-7 | Xylenes, Total | 0.56 | U | 10 | 0.56 |
| 98-06-6 | tert-Butylbenzene | 0.69 | U | 5.2 | 0.69 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP | Lab Sample ID: 240-57930-11

Matrix: Solid | Lab File ID: 146562.D

Analysis Method: 8260C | Date Collected: 11/13/2015 00:00

Sample wt/vol: 5.73(g) | Date Analyzed: 11/18/2015 21:59

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.55 | U | 5.2 | 0.55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 90 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 89 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 83 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: TB-111315 | Lab Sample ID: 240-57930-12

Matrix: Water | Lab File ID: UXX3077.D

Analysis Method: 8260C | Date Collected: 11/13/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 14:51

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.28 | U | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.44 | U | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.22 | U | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.30 | U | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 0.45 | U | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.37 | U * | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.44 | U | 1.0 | 0.44 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.47 | U | 5.0 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.32 | U | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.41 | U | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.82 | U | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.25 | U | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 0.23 | U | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 0.25 | U | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.19 | U | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 0.19 | U | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.27 | U | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 0.26 | U | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 0.53 | U | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.59 | U | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 0.40 | U | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 0.48 | U | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 10 | 0.99 |
| 67-64-1 | Acetone | 0.94 | U | 10 | 0.94 |
| 71-43-2 | Benzene | 0.35 | U | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 0.35 | U | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 0.56 | U | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 0.44 | U | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.38 | U | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 0.43 | U | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 0.25 | U | 1.0 | 0.25 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: TB-111315 | Lab Sample ID: 240-57930-12

Matrix: Water | Lab File ID: UXX3077.D

Analysis Method: 8260C | Date Collected: 11/13/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 14:51

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 1.0 | 0.43 |
| 75-00-3 | Chloroethane | 0.32 | U | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 0.25 | U | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 0.44 | U | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.26 | U | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.46 | U | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 0.42 | U | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.50 | U | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 4.9 | U | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 0.25 | U | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 0.35 | U | 1.0 | 0.35 |
| 108-20-3 | Diisopropyl ether | 0.50 | U | 10 | 0.50 |
| 98-82-8 | Isopropylbenzene | 0.35 | U | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 0.20 | U | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 0.33 | U | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.24 | U | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 0.45 | U | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 0.31 | U | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 0.40 | U | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 0.25 | U | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 0.48 | U | 1.0 | 0.48 |
| 100-42-5 | Styrene | 0.45 | U | 1.0 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.30 | U | 5.0 | 0.30 |
| 637-92-3 | Ethyl t-butyl ether | 0.23 | U | 5.0 | 0.23 |
| 127-18-4 | Tetrachloroethene | 0.31 | U | 1.0 | 0.31 |
| 108-88-3 | Toluene | 0.23 | U | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.30 | U | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.56 | U | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 0.22 | U | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 0.49 | U | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 0.41 | U | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 0.29 | U | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 0.52 | U | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 0.41 | U | 1.0 | 0.41 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: TB-111315 | Lab Sample ID: 240-57930-12

Matrix: Water | Lab File ID: UXX3077.D

Analysis Method: 8260C | Date Collected: 11/13/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 14:51

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.32 | U | 1.0 | 0.32 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 82 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 81 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 75 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 83 | | 78-125 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-206893/8 | UXX3045.D |
| Level 2 | STD8260 240-206893/7 | UXX3044.D |
| Level 3 | STD8260 240-206893/6 | UXX3043.D |
| Level 4 | STD8260 240-206893/5 | UXX3042.D |
| Level 5 | STD8260 240-206893/4 | UXX3041.D |
| Level 6 | STD8260 240-206893/3 | UXX3040.D |
| Level 7 | STD8260 240-206893/2 | UXX3039.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | 0.2558<br>0.2635 | 0.2392<br>0.2172 | 0.2617 | 0.2575 | 0.2486 | Ave | | 0.2491 | | | 0.1000 | 6.5 | | 15.0 | | | |
| Chloromethane | 0.3846<br>0.2660 | 0.3277<br>0.2465 | 0.3315 | 0.3019 | 0.2932 | Ave | | 0.3073 | | | 0.1000 | 14.9 | | 15.0 | | | |
| Vinyl chloride | 0.3532<br>0.2771 | 0.3108<br>0.2446 | 0.3169 | 0.3035 | 0.2896 | Ave | | 0.2994 | | | 0.1000 | 11.4 | | 15.0 | | | |
| Butadiene | 0.3861<br>0.2850 | 0.3426<br>0.2497 | 0.3114 | 0.3090 | 0.2873 | Ave | | 0.3102 | | | | 14.2 | | 15.0 | | | |
| Bromomethane | 0.2835<br>0.2345 | 0.2479<br>0.2522 | 0.2576 | 0.2438 | 0.2292 | Ave | | 0.2498 | | | 0.0500 | 7.1 | | 15.0 | | | |
| Chloroethane | 0.2724<br>0.2278 | 0.2797<br>0.2289 | 0.2581 | 0.2474 | 0.2415 | Ave | | 0.2508 | | | | 8.1 | | 15.0 | | | |
| Dichlorofluoromethane | 0.6581<br>0.5225 | 0.6427<br>0.5307 | 0.6029 | 0.5780 | 0.5451 | Ave | | 0.5829 | | | | 9.3 | | 15.0 | | | |
| Trichlorofluoromethane | 0.2696<br>0.3201 | 0.2402<br>0.3114 | 0.2625 | 0.2904 | 0.2679 | Ave | | 0.2803 | | | 0.1000 | 10.1 | | 15.0 | | | |
| Ethyl ether | 0.2822<br>0.2083 | 0.2470<br>0.2155 | 0.2248 | 0.2260 | 0.2246 | Ave | | 0.2326 | | | | 10.7 | | 15.0 | | | |
| Acrolein | 0.0252<br>0.0211 | 0.0281<br>0.0216 | 0.0242 | 0.0238 | 0.0231 | Ave | | 0.0238 | | | | 9.9 | | 15.0 | | | |
| 1,1-Dichloroethene | 0.2277<br>0.2160 | 0.2273<br>0.2098 | 0.2270 | 0.2229 | 0.2254 | Ave | | 0.2223 | | | 0.1000 | 3.1 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.1842<br>0.1816 | 0.1408<br>0.1610 | 0.1611 | 0.1636 | 0.1681 | Ave | | 0.1658 | | | 0.0500 | 8.8 | | 15.0 | | | |
| Acetone | 0.2427<br>0.0612 | 0.1441<br>0.0584 | 0.1048 | 0.0768 | 0.0676 | Lin1 | 0.1853 | 0.0567 | | | 0.0500 | | | | 1.0000 | | 0.9900 |
| Iodomethane | 0.3917<br>0.3480 | 0.3798<br>0.3570 | 0.3850 | 0.3783 | 0.3797 | Ave | | 0.3742 | | | | 4.2 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon disulfide | 0.6648<br>0.6777 | 0.6334<br>0.6852 | 0.6385 | 0.6682 | 0.7000 | Ave | | 0.6668 | | | 0.1000 | 3.6 | | 15.0 | | | |
| 3-Chloro-1-propene | 0.1452<br>0.1539 | 0.1544<br>0.1554 | 0.1515 | 0.1537 | 0.1637 | Ave | | 0.1540 | | | | 3.6 | | 15.0 | | | |
| Methyl acetate | 0.2076<br>0.1682 | 0.1855<br>0.1769 | 0.1829 | 0.1723 | 0.1738 | Ave | | 0.1810 | | | 0.1000 | 7.3 | | 15.0 | | | |
| Methylene Chloride | 0.8173<br>0.2350 | 0.5325<br>0.2430 | 0.3803 | 0.2969 | 0.2707 | Lin1 | 0.2967 | 0.2323 | | | 0.1000 | | | | 0.9990 | | 0.9900 |
| tert-Butyl alcohol (TBA) | 0.0163<br>0.0168 | 0.0190<br>0.0156 | 0.0171 | 0.0180 | 0.0176 | Ave | | 0.0172 | | | | 6.5 | | 15.0 | | | |
| Acrylonitrile | 0.0901<br>0.0885 | 0.0917<br>0.0936 | 0.0885 | 0.0895 | 0.0897 | Ave | | 0.0902 | | | | 2.0 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.7089<br>0.7125 | 0.6937<br>0.7861 | 0.6936 | 0.7027 | 0.7319 | Ave | | 0.7185 | | | 0.1000 | 4.5 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.2637<br>0.2513 | 0.2813<br>0.2576 | 0.2574 | 0.2537 | 0.2619 | Ave | | 0.2610 | | | 0.1000 | 3.8 | | 15.0 | | | |
| Hexane | 0.0667<br>0.0591 | 0.0445<br>0.0515 | 0.0468 | 0.0549 | 0.0504 | Ave | | 0.0534 | | | | 14.2 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.4688<br>0.4268 | 0.4868<br>0.4444 | 0.4594 | 0.4569 | 0.4642 | Ave | | 0.4582 | | | 0.1000 | 4.1 | | 15.0 | | | |
| Vinyl acetate | 0.4652<br>0.4501 | 0.4809<br>0.4985 | 0.4063 | 0.4376 | 0.4468 | Ave | | 0.4551 | | | | 6.6 | | 15.0 | | | |
| 2,2-Dichloropropane | 0.2507<br>0.2589 | 0.2505<br>0.2597 | 0.2625 | 0.2646 | 0.2765 | Ave | | 0.2605 | | | | 3.4 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.2875<br>0.2595 | 0.3082<br>0.2632 | 0.2741 | 0.2812 | 0.2787 | Ave | | 0.2789 | | | 0.1000 | 5.8 | | 15.0 | | | |
| 2-Butanone (MEK) | 0.1329<br>0.0882 | 0.1145<br>0.0909 | 0.1044 | 0.0968 | 0.0950 | Lin1 | 0.0483 | 0.0899 | | | 0.0500 | | | | 0.9990 | | 0.9900 |
| Bromochloromethane | 0.1557<br>0.1254 | 0.1482<br>0.1269 | 0.1350 | 0.1410 | 0.1396 | Ave | | 0.1388 | | | | 7.9 | | 15.0 | | | |
| Tetrahydrofuran | 0.0810<br>0.0635 | 0.0739<br>0.0622 | 0.0622 | 0.0658 | 0.0659 | Ave | | 0.0678 | | | | 10.4 | | 15.0 | | | |
| Chloroform | 0.4458<br>0.4173 | 0.4612<br>0.4267 | 0.4424 | 0.4440 | 0.4443 | Ave | | 0.4402 | | | | 3.2 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.3394<br>0.3314 | 0.3153<br>0.3247 | 0.3147 | 0.3438 | 0.3454 | Ave | | 0.3307 | | | 0.1000 | 3.9 | | 15.0 | | | |
| Cyclohexane | 0.3340<br>0.3898 | 0.2912<br>0.3592 | 0.3583 | 0.3638 | 0.3749 | Ave | | 0.3530 | | | 0.1000 | 9.1 | | 15.0 | | | |
| 1,1-Dichloropropene | 0.3273<br>0.3093 | 0.3070<br>0.3044 | 0.3128 | 0.3083 | 0.3137 | Ave | | 0.3118 | | | | 2.4 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon tetrachloride | 0.2581<br>0.2939 | 0.2554<br>0.2921 | 0.2573 | 0.2680 | 0.2909 | Ave | | 0.2737 | | | 0.1000 | 6.5 | | 15.0 | | | |
| Isobutyl alcohol | 0.0106<br>0.0096 | 0.0118<br>0.0090 | 0.0096 | 0.0092 | 0.0096 | Ave | | 0.0099 | | | | 9.8 | | 15.0 | | | |
| Benzene | 1.1454<br>0.9777 | 1.0652<br>1.0667 | 1.0301 | 1.0196 | 1.0518 | Ave | | 1.0509 | | | 0.5000 | 4.9 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.3885<br>0.3324 | 0.3590<br>0.3530 | 0.3467 | 0.3486 | 0.3519 | Ave | | 0.3543 | | | 0.1000 | 4.8 | | 15.0 | | | |
| n-Heptane | 0.0493<br>0.0489 | 0.0294<br>0.0431 | 0.0386 | 0.0414 | 0.0396 | Lin1 | -0.005 | 0.0442 | | | | | | | 0.9930 | | 0.9900 |
| Trichloroethene | 0.2751<br>0.2556 | 0.2617<br>0.2592 | 0.2722 | 0.2658 | 0.2730 | Ave | | 0.2661 | | | 0.2000 | 2.8 | | 15.0 | | | |
| Methylcyclohexane | 0.3025<br>0.3628 | 0.2590<br>0.3315 | 0.2976 | 0.3223 | 0.3257 | Ave | | 0.3145 | | | 0.1000 | 10.3 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.2503<br>0.2144 | 0.2422<br>0.2285 | 0.2313 | 0.2318 | 0.2407 | Ave | | 0.2342 | | | 0.1000 | 4.9 | | 15.0 | | | |
| Dibromomethane | 0.1686<br>0.1359 | 0.1537<br>0.1431 | 0.1427 | 0.1473 | 0.1491 | Ave | | 0.1487 | | | | 7.0 | | 15.0 | | | |
| 1,4-Dioxane | 0.0027<br>0.0024 | 0.0023<br>0.0021 | 0.0025 | 0.0027 | 0.0027 | Ave | | 0.0025 | | | | 9.7 | | 15.0 | | | |
| Bromodichloromethane | 0.2753<br>0.2946 | 0.2940<br>0.3226 | 0.2839 | 0.2895 | 0.3147 | Ave | | 0.2964 | | | 0.1000 | 5.6 | | 15.0 | | | |
| 2-Chloroethyl vinyl ether | 0.1105<br>0.1122 | 0.1074<br>0.1258 | 0.1121 | 0.1153 | 0.1234 | Ave | | 0.1152 | | | | 6.0 | | 15.0 | | | |
| cis-1,3-Dichloropropene | 0.2903<br>0.3349 | 0.2953<br>0.3723 | 0.3095 | 0.3263 | 0.3672 | Ave | | 0.3280 | | | 0.1500 | 9.9 | | 15.0 | | | |
| 4-Methyl-2-pentanone (MIBK) | 0.1680<br>0.1918 | 0.1902<br>0.1990 | 0.1896 | 0.1873 | 0.2030 | Ave | | 0.1898 | | | 0.0500 | 5.9 | | 15.0 | | | |
| Toluene | 1.5017<br>1.4446 | 1.4781<br>1.6471 | 1.4250 | 1.4205 | 1.5133 | Ave | | 1.4900 | | | 0.4000 | 5.2 | | 15.0 | | | |
| trans-1,3-Dichloropropene | 0.3488<br>0.4112 | 0.3467<br>0.4578 | 0.3525 | 0.3902 | 0.4305 | Ave | | 0.3911 | | | 0.1000 | 11.3 | | 15.0 | | | |
| Ethyl methacrylate | 0.2662<br>0.3456 | 0.2655<br>0.3743 | 0.2760 | 0.3068 | 0.3532 | Ave | | 0.3125 | | | | 14.5 | | 15.0 | | | |
| 1,1,2-Trichloroethane | 0.3131<br>0.2615 | 0.2797<br>0.2709 | 0.2736 | 0.2823 | 0.2862 | Ave | | 0.2810 | | | 0.1000 | 5.8 | | 15.0 | | | |
| Tetrachloroethene | 0.3330<br>0.2841 | 0.2857<br>0.2885 | 0.2818 | 0.2781 | 0.2967 | Ave | | 0.2926 | | | 0.2000 | 6.4 | | 15.0 | | | |
| 1,3-Dichloropropane | 0.5246<br>0.4664 | 0.5114<br>0.4971 | 0.5059 | 0.4948 | 0.5021 | Ave | | 0.5003 | | | | 3.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2-Hexanone | 0.1672<br>0.1736 | 0.1699<br>0.1723 | 0.1603 | 0.1688 | 0.1761 | Ave | | 0.1698 | | | 0.0500 | 3.0 | | 15.0 | | | |
| Dibromochloromethane | 0.2366<br>0.2993 | 0.2567<br>0.3306 | 0.2661 | 0.2732 | 0.3096 | Ave | | 0.2817 | | | | 11.6 | | 15.0 | | | |
| 1,2-Dibromoethane | 0.2764<br>0.2724 | 0.2646<br>0.2837 | 0.2503 | 0.2739 | 0.2865 | Ave | | 0.2725 | | | | 4.5 | | 15.0 | | | |
| Chlorobenzene | 0.9905<br>0.9215 | 0.9541<br>0.9901 | 0.9508 | 0.9618 | 0.9770 | Ave | | 0.9637 | | | 0.3000 | 2.6 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.2776<br>0.3303 | 0.2881<br>0.3559 | 0.3154 | 0.3218 | 0.3460 | Ave | | 0.3193 | | | | 9.0 | | 15.0 | | | |
| Ethylbenzene | 0.4968<br>0.5027 | 0.4828<br>0.5356 | 0.4808 | 0.4888 | 0.5305 | Ave | | 0.5026 | | | | 4.4 | | 15.0 | | | |
| m-Xylene & p-Xylene | 0.5798<br>0.6297 | 0.5575<br>0.6639 | 0.5971 | 0.6068 | 0.5733 | Ave | | 0.6012 | | | | 6.0 | | 15.0 | | | |
| o-Xylene | 0.5320<br>0.6217 | 0.5651<br>0.6762 | 0.5750 | 0.6049 | 0.6036 | Ave | | 0.5969 | | | | 7.7 | | 15.0 | | | |
| Styrene | 0.8405<br>1.0572 | 0.8487<br>1.1974 | 0.8768 | 0.9758 | 1.0157 | Ave | | 0.9732 | | | 0.3000 | 13.4 | | 15.0 | | | |
| Bromoform | 0.1571<br>0.2004 | 0.1575<br>0.2291 | 0.1631 | 0.1736 | 0.1945 | Ave | | 0.1822 | | | 0.1000 | 14.8 | | 15.0 | | | |
| Isopropylbenzene | 1.3238<br>1.6576 | 1.2674<br>1.8024 | 1.4275 | 1.5101 | 1.5793 | Ave | | 1.5097 | | | 0.1000 | 12.5 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 0.6360<br>0.6040 | 0.6329<br>0.5976 | 0.6209 | 0.6110 | 0.6383 | Ave | | 0.6201 | | | 0.3000 | 2.6 | | 15.0 | | | |
| Bromobenzene | 0.8001<br>0.6932 | 0.7459<br>0.7093 | 0.7303 | 0.7276 | 0.7324 | Ave | | 0.7341 | | | | 4.6 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.2183<br>0.2038 | 0.2576<br>0.1962 | 0.2137 | 0.2194 | 0.2171 | Ave | | 0.2180 | | | | 8.9 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | 0.2075<br>0.1998 | 0.1781<br>0.1966 | 0.1830 | 0.1661 | 0.2009 | Ave | | 0.1903 | | | | 7.8 | | 15.0 | | | |
| N-Propylbenzene | 0.6116<br>0.7284 | 0.6386<br>0.7506 | 0.6933 | 0.7402 | 0.7626 | Ave | | 0.7036 | | | | 8.3 | | 15.0 | | | |
| 2-Chlorotoluene | 0.6566<br>0.6365 | 0.6780<br>0.6434 | 0.6623 | 0.6364 | 0.6860 | Ave | | 0.6570 | | | | 3.0 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 1.6963<br>2.1622 | 1.8455<br>2.3447 | 1.9011 | 2.0426 | 2.2341 | Ave | | 2.0324 | | | | 11.4 | | 15.0 | | | |
| 4-Chlorotoluene | 0.6439<br>0.6935 | 0.6990<br>0.7194 | 0.6656 | 0.6925 | 0.7421 | Ave | | 0.6937 | | | | 4.7 | | 15.0 | | | |
| tert-Butylbenzene | 1.6145<br>1.8645 | 1.5963<br>1.9300 | 1.7599 | 1.7724 | 1.9262 | Ave | | 1.7806 | | | | 7.7 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,2,4-Trimethylbenzene | 2.1077<br>2.2890 | 2.0342<br>2.4728 | 2.0415 | 2.1970 | 2.4160 | Ave | | 2.2226 | | | | 7.9 | | 15.0 | | | |
| sec-Butylbenzene | 1.9388<br>2.4846 | 2.1534<br>2.6180 | 2.1867 | 2.3241 | 2.5155 | Ave | | 2.3173 | | | | 10.3 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.4675<br>1.3090 | 1.3805<br>1.3608 | 1.3510 | 1.3428 | 1.4138 | Ave | | 1.3751 | | | 0.6000 | 3.8 | | 15.0 | | | |
| p-Isopropyltoluene | 1.7191<br>2.2370 | 1.7806<br>2.4373 | 1.9795 | 2.0723 | 2.2708 | Ave | | 2.0709 | | | | 12.7 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 1.5021<br>1.3812 | 1.5177<br>1.4129 | 1.4182 | 1.4017 | 1.4978 | Ave | | 1.4474 | | | 0.5000 | 3.9 | | 15.0 | | | |
| n-Butylbenzene | 1.3933<br>1.7156 | 1.4227<br>1.7988 | 1.4703 | 1.6110 | 1.7497 | Ave | | 1.5945 | | | | 10.4 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.3196<br>1.2764 | 1.3719<br>1.2867 | 1.3402 | 1.3154 | 1.3934 | Ave | | 1.3291 | | | 0.4000 | 3.2 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | 0.0961<br>0.1330 | 0.0948<br>0.1119 | 0.1082 | 0.1009 | 0.1141 | Ave | | 0.1084 | | | 0.0500 | 12.1 | | 15.0 | | | |
| 1,2,4-Trichlorobenzene | 0.8976<br>0.7771 | 0.8558<br>0.7077 | 0.8693 | 0.8206 | 0.8326 | Ave | | 0.8230 | | | 0.2000 | 7.7 | | 15.0 | | | |
| Hexachlorobutadiene | 0.3143<br>0.2739 | 0.3028<br>0.2463 | 0.3025 | 0.2992 | 0.2993 | Ave | | 0.2912 | | | | 8.0 | | 15.0 | | | |
| Naphthalene | 1.8805<br>2.1523 | 2.0421<br>1.8342 | 2.0598 | 2.0535 | 2.1417 | Ave | | 2.0234 | | | | 6.0 | | 15.0 | | | |
| 1,2,3-Trichlorobenzene | 0.8427<br>0.7352 | 0.8414<br>0.6087 | 0.8664 | 0.7993 | 0.8150 | Ave | | 0.7870 | | | | 11.3 | | 15.0 | | | |
| Dibromofluoromethane (Surr) | 0.2463<br>0.2220 | 0.2463<br>0.2239 | 0.2339 | 0.2284 | 0.2375 | Ave | | 0.2341 | | | | 4.2 | | 15.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | 0.3305<br>0.2719 | 0.3160<br>0.2737 | 0.2937 | 0.2907 | 0.2945 | Ave | | 0.2958 | | | | 7.2 | | 15.0 | | | |
| Toluene-d8 (Surr) | 1.3377<br>1.2414 | 1.2237<br>1.3339 | 1.1875 | 1.2376 | 1.2927 | Ave | | 1.2650 | | | | 4.5 | | 15.0 | | | |
| 4-Bromofluorobenzene (Surr) | 0.4225<br>0.4246 | 0.4122<br>0.4347 | 0.4494 | 0.4250 | 0.4401 | Ave | | 0.4298 | | | | 2.9 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-206893/8 | UXX3045.D |
| Level 2 | STD8260 240-206893/7 | UXX3044.D |
| Level 3 | STD8260 240-206893/6 | UXX3043.D |
| Level 4 | STD8260 240-206893/5 | UXX3042.D |
| Level 5 | STD8260 240-206893/4 | UXX3041.D |
| Level 6 | STD8260 240-206893/3 | UXX3040.D |
| Level 7 | STD8260 240-206893/2 | UXX3039.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 |
| Dichlorodifluoromethane | FB | Ave | 17265<br>898833 | 33006<br>1600119 | 74869 | 191559 | 391222 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Chloromethane | FB | Ave | 25953<br>907247 | 45204<br>1815442 | 94867 | 224619 | 461389 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Vinyl chloride | FB | Ave | 23834<br>945047 | 42885<br>1801704 | 90673 | 225779 | 455798 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Butadiene | FB | Ave | 26056<br>972071 | 47271<br>1839739 | 89091 | 229854 | 452088 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Bromomethane | FB | Ave | 19132<br>799715 | 34204<br>1857581 | 73707 | 181401 | 360681 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Chloroethane | FB | Ave | 18385<br>776852 | 38584<br>1686308 | 73842 | 184077 | 380132 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Dichlorofluoromethane | FB | Ave | 44411<br>1782274 | 88674<br>3909191 | 172515 | 430010 | 857882 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Trichlorofluoromethane | FB | Ave | 18196<br>1091877 | 33141<br>2293874 | 75101 | 216066 | 421680 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Ethyl ether | FB | Ave | 19045<br>710563 | 34078<br>1587337 | 64338 | 168130 | 353478 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Acrolein | FB | Ave | 8489<br>359160 | 19376<br>795558 | 34588 | 88407 | 181579 | 2.50<br>100 | 5.00<br>200 | 10.0 | 25.0 | 50.0 |
| 1,1-Dichloroethene | FB | Ave | 15365<br>736872 | 31365<br>1545523 | 64962 | 165790 | 354730 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroetha ne | FB | Ave | 12434<br>619515 | 19429<br>1185805 | 46085 | 121672 | 264579 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Acetone | FB | Lin1 | 32759<br>417420 | 39766<br>861052 | 59957 | 114218 | 212684 | 1.00<br>40.0 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 |
| Iodomethane | FB | Ave | 26435<br>1186825 | 52396<br>2629420 | 110159 | 281413 | 597660 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Carbon disulfide | FB | Ave | 44863<br>2311463 | 87391<br>5047429 | 182691 | 497083 | 1101772 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 |
| 3-Chloro-1-propene | FB | Ave | 9797<br>524913 | 21296<br>1144481 | 43357 | 114365 | 257638 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Methyl acetate | FB | Ave | 70054<br>2868384 | 127982<br>6513770 | 261663 | 640721 | 1367770 | 2.50<br>100 | 5.00<br>200 | 10.0 | 25.0 | 50.0 |
| Methylene Chloride | FB | Lin1 | 55155<br>801591 | 73466<br>1789870 | 108826 | 220889 | 426067 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| tert-Butyl alcohol (TBA) | FB | Ave | 10983<br>574602 | 26180<br>1151497 | 49070 | 134137 | 276917 | 5.00<br>200 | 10.0<br>400 | 20.0 | 50.0 | 100 |
| Acrylonitrile | FB | Ave | 60783<br>3020204 | 126451<br>6897566 | 253134 | 665910 | 1411077 | 5.00<br>200 | 10.0<br>400 | 20.0 | 50.0 | 100 |
| Methyl tert-butyl ether | FB | Ave | 47842<br>2430270 | 95701<br>5790520 | 198467 | 522762 | 1151925 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| trans-1,2-Dichloroethene | FB | Ave | 17799<br>857269 | 38806<br>1897485 | 73658 | 188763 | 412242 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Hexane | FB | Ave | 4500<br>201566 | 6140<br>379634 | 13398 | 40819 | 79306 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,1-Dichloroethane | FB | Ave | 31639<br>1455854 | 67157<br>3273555 | 131442 | 339908 | 730534 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Vinyl acetate | FB | Ave | 31398<br>1535147 | 66340<br>3672265 | 116260 | 325558 | 703234 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 2,2-Dichloropropane | FB | Ave | 16919<br>882981 | 34553<br>1912854 | 75111 | 196812 | 435212 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| cis-1,2-Dichloroethene | FB | Ave | 19401<br>884971 | 42526<br>1938749 | 78432 | 209224 | 438568 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 2-Butanone (MEK) | FB | Lin1 | 17932<br>601483 | 31582<br>1339128 | 59718 | 143982 | 298909 | 1.00<br>40.0 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 |
| Bromochloromethane | FB | Ave | 10507<br>427699 | 20452<br>934969 | 38634 | 104868 | 219653 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Tetrahydrofuran | FB | Ave | 10930<br>433023 | 20404<br>916582 | 35621 | 97826 | 207524 | 1.00<br>40.0 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 |
| Chloroform | FB | Ave | 30089<br>1423153 | 63632<br>3143103 | 126576 | 330276 | 699312 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,1,1-Trichloroethane | FB | Ave | 22908<br>1130178 | 43503<br>2391728 | 90053 | 255755 | 543557 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Cyclohexane | FB | Ave | 22544<br>1329454 | 40179<br>2645663 | 102526 | 270634 | 590080 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,1-Dichloropropene | FB | Ave | 22086<br>1054925 | 42352<br>2242384 | 89507 | 229387 | 493783 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Carbon tetrachloride | FB | Ave | 17421<br>1002356 | 35236<br>2151640 | 73629 | 199382 | 457894 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Isobutyl alcohol | CBZ | Ave | 12612<br>569173 | 29509<br>1174558 | 49336 | 121468 | 270636 | 12.5<br>500 | 25.0<br>1000 | 50.0 | 125 | 250 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 |
| Benzene | FB | Ave | 77298<br>3334803 | 146955<br>7857895 | 294765 | 758516 | 1655327 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2-Dichloroethane | FB | Ave | 26217<br>1133832 | 49525<br>2600255 | 99195 | 259306 | 553764 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| n-Heptane | FB | Lin1 | 3325<br>166762 | 4058<br>317276 | 11032 | 30785 | 62361 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Trichloroethene | FB | Ave | 18563<br>871874 | 36102<br>1909675 | 77883 | 197763 | 429652 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Methylcyclohexane | FB | Ave | 20417<br>1237549 | 35733<br>2441887 | 85142 | 239788 | 512564 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2-Dichloropropane | FB | Ave | 16894<br>731313 | 33414<br>1683412 | 66182 | 172428 | 378861 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Dibromomethane | FB | Ave | 11381<br>463530 | 21210<br>1054246 | 40844 | 109609 | 234626 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,4-Dioxane | FB | Ave | 3602<br>166992 | 6320<br>302882 | 14382 | 40077 | 84876 | 10.0<br>400 | 20.0<br>800 | 40.0 | 100 | 200 |
| Bromodichloromethane | FB | Ave | 18579<br>1004770 | 40564<br>2376308 | 81222 | 215341 | 495322 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 2-Chloroethyl vinyl ether | FB | Ave | 14919<br>765094 | 29621<br>1853882 | 64179 | 171500 | 388433 | 1.00<br>40.0 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 |
| cis-1,3-Dichloropropene | FB | Ave | 19595<br>1142246 | 40738<br>2742359 | 88563 | 242733 | 577839 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 4-Methyl-2-pentanone (MIBK) | FB | Ave | 22677<br>1308361 | 52471<br>2931616 | 108492 | 278647 | 639088 | 1.00<br>40.0 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 |
| Toluene | CBZ | Ave | 71518<br>3417455 | 147451<br>8551971 | 293542 | 750449 | 1705813 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| trans-1,3-Dichloropropene | CBZ | Ave | 16613<br>972870 | 34584<br>2376876 | 72622 | 206130 | 485276 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Ethyl methacrylate | CBZ | Ave | 12678<br>817653 | 26483<br>1943325 | 56850 | 162055 | 398125 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,1,2-Trichloroethane | CBZ | Ave | 14909<br>618556 | 27906<br>1406801 | 56358 | 149135 | 322593 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Tetrachloroethene | CBZ | Ave | 15858<br>672151 | 28500<br>1498132 | 58058 | 146922 | 334467 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,3-Dichloropropane | CBZ | Ave | 24983<br>1103440 | 51017<br>2580934 | 104220 | 261387 | 565946 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 2-Hexanone | CBZ | Ave | 15929<br>821590 | 33891<br>1789608 | 66054 | 178392 | 397090 | 1.00<br>40.0 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 |
| Dibromochloromethane | CBZ | Ave | 11266<br>708035 | 25612<br>1716296 | 54814 | 144317 | 349045 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2-Dibromoethane | CBZ | Ave | 13165<br>644368 | 26396<br>1472985 | 51567 | 144697 | 322932 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 |
| Chlorobenzene | CBZ | Ave | 47172<br>2180031 | 95179<br>5140823 | 195854 | 508109 | 1101321 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,1,1,2-Tetrachloroethane | CBZ | Ave | 13219<br>781292 | 28739<br>1847738 | 64970 | 170005 | 390016 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Ethylbenzene | CBZ | Ave | 23661<br>1189183 | 48166<br>2781160 | 99045 | 258233 | 598009 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| m-Xylene & p-Xylene | CBZ | Ave | 27613<br>1489611 | 55612<br>3447096 | 122998 | 320538 | 646301 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| o-Xylene | CBZ | Ave | 25337<br>1470848 | 56368<br>3510865 | 118451 | 319583 | 680416 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Styrene | CBZ | Ave | 40028<br>2501041 | 84665<br>6217212 | 180628 | 515514 | 1144967 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Bromoform | CBZ | Ave | 7483<br>474080 | 15707<br>1189437 | 33595 | 91697 | 219244 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Isopropylbenzene | CBZ | Ave | 63042<br>3921294 | 126433<br>9358353 | 294064 | 797791 | 1780215 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,1,2,2-Tetrachloroethane | DCB | Ave | 17183<br>873113 | 35340<br>1921331 | 73647 | 193446 | 423089 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Bromobenzene | DCB | Ave | 21616<br>1002060 | 41651<br>2280655 | 86623 | 230348 | 485443 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2,3-Trichloropropane | DCB | Ave | 5898<br>294541 | 14385<br>630912 | 25345 | 69463 | 143889 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| trans-1,4-Dichloro-2-butene | DCB | Ave | 5606<br>288879 | 9945<br>632232 | 21702 | 52571 | 133183 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| N-Propylbenzene | DCB | Ave | 16522<br>1052947 | 35662<br>2413401 | 82228 | 234349 | 505489 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 2-Chlorotoluene | DCB | Ave | 17738<br>920064 | 37858<br>2068657 | 78556 | 201473 | 454727 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,3,5-Trimethylbenzene | DCB | Ave | 45828<br>3125380 | 103057<br>7538766 | 225491 | 646650 | 1480857 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 4-Chlorotoluene | DCB | Ave | 17395<br>1002418 | 39034<br>2313127 | 78948 | 219229 | 491892 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| tert-Butylbenzene | DCB | Ave | 43619<br>2695135 | 89141<br>6205575 | 208741 | 561094 | 1276733 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2,4-Trimethylbenzene | DCB | Ave | 56942<br>3308757 | 113592<br>7950697 | 242142 | 695541 | 1601393 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| sec-Butylbenzene | DCB | Ave | 52380<br>3591513 | 120249<br>8417653 | 259369 | 735761 | 1667345 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,3-Dichlorobenzene | DCB | Ave | 39647<br>1892165 | 77090<br>4375383 | 160243 | 425101 | 937107 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| p-Isopropyltoluene | DCB | Ave | 46443<br>3233508 | 99428<br>7836422 | 234790 | 656047 | 1505160 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 09:23 Calibration End Date: 11/16/2015 11:35 Calibration ID: 31566

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3 | LVL 4 | LVL 5 |
| 1,4-Dichlorobenzene | DCB | Ave | 40581<br>1996520 | 84749<br>4542939 | 168209 | 443754 | 992814 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| n-Butylbenzene | DCB | Ave | 37643<br>2479891 | 79444<br>5783584 | 174386 | 510006 | 1159746 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2-Dichlorobenzene | DCB | Ave | 35650<br>1845033 | 76608<br>4137023 | 158959 | 416422 | 923596 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2-Dibromo-3-Chloropropane | DCB | Ave | 2597<br>192198 | 5294<br>359844 | 12835 | 31935 | 75606 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2,4-Trichlorobenzene | DCB | Ave | 24251<br>1123256 | 47787<br>2275298 | 103111 | 259789 | 551862 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Hexachlorobutadiene | DCB | Ave | 8491<br>395958 | 16907<br>791778 | 35885 | 94721 | 198391 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Naphthalene | DCB | Ave | 50804<br>3111111 | 114032<br>5897375 | 244314 | 650107 | 1419614 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2,3-Trichlorobenzene | DCB | Ave | 22766<br>1062726 | 46984<br>1957067 | 102768 | 253046 | 540232 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Dibromofluoromethane (Surr) | FB | Ave | 16622<br>757265 | 33976<br>1649233 | 66937 | 169930 | 373851 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Ave | 22304<br>927283 | 43593<br>2016090 | 84041 | 216265 | 463511 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| Toluene-d8 (Surr) | CBZ | Ave | 63707<br>2936784 | 122072<br>6926103 | 244634 | 653827 | 1457157 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |
| 4-Bromofluorobenzene (Surr) | CBZ | Ave | 20122<br>1004511 | 41124<br>2256868 | 92570 | 224511 | 496055 | 0.500<br>20.0 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin1 = Linear 1/conc ISTD |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 12:19 Calibration End Date: 11/16/2015 14:10 Calibration ID: 31568

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STDA9 240-206893/15 | UXX3052.D |
| Level 2 | STDA9 240-206893/14 | UXX3051.D |
| Level 3 | STDA9 240-206893/13 | UXX3050.D |
| Level 4 | STDA9 240-206893/12 | UXX3049.D |
| Level 5 | STDA9 240-206893/11 | UXX3048.D |
| Level 6 | STDA9 240-206893/10 | UXX3047.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Acetonitrile | 0.0271 0.0239 | 0.0277 | 0.0237 | 0.0273 | 0.0243 | Ave | | 0.0257 | | | | 7.3 | | 15.0 | | | |
| Diisopropyl ether | 0.2074 0.2141 | 0.1954 | 0.2019 | 0.2122 | 0.2041 | Ave | | 0.2058 | | | | 3.4 | | 15.0 | | | |
| 2-Chloro-1,3-butadiene | 0.3039 0.3467 | 0.2883 | 0.3203 | 0.3290 | 0.3231 | Ave | | 0.3186 | | | | 6.4 | | 15.0 | | | |
| Ethyl t-butyl ether | 0.6075 0.7078 | 0.5728 | 0.6237 | 0.6532 | 0.6576 | Ave | | 0.6371 | | | | 7.3 | | 15.0 | | | |
| Ethyl acetate | 0.2145 0.1898 | 0.2036 | 0.1879 | 0.1826 | 0.1831 | Ave | | 0.1936 | | | | 6.6 | | 15.0 | | | |
| Propionitrile | 0.0331 0.0315 | 0.0310 | 0.0322 | 0.0316 | 0.0313 | Ave | | 0.0318 | | | | 2.4 | | 15.0 | | | |
| Methacrylonitrile | 0.1369 0.1768 | 0.1321 | 0.1400 | 0.1443 | 0.1510 | Ave | | 0.1468 | | | | 10.9 | | 15.0 | | | |
| Tert-amyl methyl ether | 0.5203 0.7076 | 0.5316 | 0.5643 | 0.6101 | 0.6307 | Ave | | 0.5941 | | | | 11.8 | | 15.0 | | | |
| n-Butanol | 0.0064 0.0067 | 0.0058 | 0.0063 | 0.0072 | 0.0068 | Ave | | 0.0065 | | | | 7.4 | | 15.0 | | | |
| Ethyl acrylate | 0.2087 0.2487 | 0.2044 | 0.2128 | 0.2199 | 0.2344 | Ave | | 0.2215 | | | | 7.7 | | 15.0 | | | |
| Methyl methacrylate | 0.1502 0.1925 | 0.1500 | 0.1715 | 0.1728 | 0.1777 | Ave | | 0.1691 | | | | 9.8 | | 15.0 | | | |
| 2-Nitropropane | 0.0327 0.0486 | 0.0380 | 0.0364 | 0.0356 | 0.0393 | Ave | | 0.0384 | | | | 14.3 | | 15.0 | | | |
| n-Butyl acetate | 0.1150 0.1544 | 0.1197 | 0.1313 | 0.1462 | 0.1490 | Ave | | 0.1359 | | | | 12.0 | | 15.0 | | | |
| 1-Chlorohexane | 0.3404 0.3832 | 0.3136 | 0.3413 | 0.3615 | 0.3608 | Ave | | 0.3501 | | | | 6.8 | | 15.0 | | | |
| Cyclohexanone | 0.0126 0.0125 | 0.0121 | 0.0129 | 0.0147 | 0.0127 | Ave | | 0.0129 | | | | 7.1 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 12:19 Calibration End Date: 11/16/2015 14:10 Calibration ID: 31568

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,2,3-Trimethylbenzene | 1.8196<br>2.6705 | 1.8734 | 1.9730 | 2.2718 | 2.3585 | Lin1 | -1.283 | 2.5620 | | | | | | | 0.9940 | | 0.9900 |
| 1,3,5-Trichlorobenzene | 0.9002<br>0.8711 | 0.8419 | 0.6795 | 0.9108 | 0.8712 | Ave | | 0.8458 | | | | 10.1 | | 15.0 | | | |
| 2-Methylnaphthalene | 0.8988<br>0.8525 | 0.9845 | 0.9275 | 1.1949 | 1.1589 | Ave | | 1.0028 | | | | 14.1 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 12:19 Calibration End Date: 11/16/2015 14:10 Calibration ID: 31568

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STDA9 240-206893/15 | UXX3052.D |
| Level 2 | STDA9 240-206893/14 | UXX3051.D |
| Level 3 | STDA9 240-206893/13 | UXX3050.D |
| Level 4 | STDA9 240-206893/12 | UXX3049.D |
| Level 5 | STDA9 240-206893/11 | UXX3048.D |
| Level 6 | STDA9 240-206893/10 | UXX3047.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Acetonitrile | FB | Ave | 33784<br>1406719 | 69938 | 154507 | 367519 | 696514 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Diisopropyl ether | FB | Ave | 25875<br>1262849 | 49353 | 131398 | 285976 | 584031 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Chloro-1,3-butadiene | FB | Ave | 37925<br>2044723 | 72816 | 208447 | 443487 | 924411 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl t-butyl ether | FB | Ave | 75813<br>4174265 | 144670 | 405956 | 880359 | 1881458 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl acetate | FB | Ave | 53521<br>2238994 | 102847 | 244538 | 492326 | 1047697 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Propionitrile | FB | Ave | 41349<br>1857261 | 78251 | 209833 | 425524 | 895778 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Methacrylonitrile | FB | Ave | 170885<br>10427349 | 333590 | 911162 | 1944419 | 4319064 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Tert-amyl methyl ether | FB | Ave | 64925<br>4173212 | 134252 | 367320 | 822282 | 1804433 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| n-Butanol | CBZ | Ave | 14529<br>717235 | 26592 | 75298 | 175373 | 343307 | 25.0<br>1000 | 50.0 | 125 | 250 | 500 |
| Ethyl acrylate | FB | Ave | 26045<br>1466405 | 51621 | 138538 | 296429 | 670569 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Methyl methacrylate | FB | Ave | 37489<br>2270268 | 75777 | 223274 | 465950 | 1016553 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| 2-Nitropropane | FB | Ave | 8160<br>573588 | 19212 | 47396 | 95896 | 225134 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| n-Butyl acetate | CBZ | Ave | 10448<br>658195 | 22007 | 62743 | 142693 | 299255 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1-Chlorohexane | CBZ | Ave | 30919<br>1633936 | 57637 | 163044 | 352854 | 724728 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Cyclohexanone | DCB | Ave | 6117<br>272895 | 11951 | 30713 | 77762 | 137248 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| 1,2,3-Trimethylbenzene | DCB | Lin1 | 88262<br>5821766 | 184612 | 471200 | 1197911 | 2539904 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206893

SDG No.:

Instrument ID: A3UX10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 12:19 Calibration End Date: 11/16/2015 14:10 Calibration ID: 31568

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| 1,3,5-Trichlorobenzene | DCB | Ave | 43663<br>1899076 | 82961 | 162288 | 480277 | 938233 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Methylnaphthalene | DCB | Ave | 87193<br>3717072 | 194029 | 443041 | 1260113 | 2496110 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin1 = Linear 1/conc ISTD |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50 Calibration End Date: 08/13/2015 20:42 Calibration ID: 30158

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-193150/20 | 144584.D |
| Level 2 | IC 240-193150/19 | 144583.D |
| Level 3 | IC 240-193150/18 | 144582.D |
| Level 4 | IC 240-193150/17 | 144581.D |
| Level 5 | IC 240-193150/16 | 144580.D |
| Level 6 | IC 240-193150/15 | 144579.D |
| Level 7 | IC 240-193150/14 | 144578.D |
| Level 8 | IC 240-193150/13 | 144577.D |
| Level 9 | IC 240-193150/12 | 144576.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Acetonitrile | 0.0270<br>0.0217 | 0.0199<br>0.0226 | 0.0223<br>0.0219 | 0.0211<br>0.0222 | 0.0231 | Ave | | 0.0224 | | | | 8.8 | | 15.0 | | | |
| 2-Chloro-1,3-butadiene | 0.3946<br>0.4330 | 0.4036<br>0.4520 | 0.4843<br>0.4367 | 0.4330<br>0.4384 | 0.4314 | Ave | | 0.4341 | | | | 6.0 | | 15.0 | | | |
| Diisopropyl ether | 0.1991<br>0.2087 | 0.1920<br>0.2170 | 0.2189<br>0.2067 | 0.2161<br>0.2066 | 0.2062 | Ave | | 0.2079 | | | | 4.2 | | 15.0 | | | |
| Ethyl t-butyl ether | 0.8061<br>0.8145 | 0.8124<br>0.8609 | 0.8459<br>0.8423 | 0.8104<br>0.8475 | 0.7961 | Ave | | 0.8262 | | | | 2.8 | | 15.0 | | | |
| Ethyl acetate | 0.2082<br>0.1555 | 0.1520<br>0.1600 | 0.1502<br>0.1572 | 0.1518<br>0.1578 | 0.1525 | Ave | | 0.1606 | | | | 11.3 | | 15.0 | | | |
| Propionitrile | 0.0291<br>0.0267 | 0.0294<br>0.0281 | 0.0277<br>0.0272 | 0.0279<br>0.0272 | 0.0264 | Ave | | 0.0278 | | | | 3.7 | | 15.0 | | | |
| Methacrylonitrile | 0.1071<br>0.1082 | 0.1056<br>0.1106 | 0.1086<br>0.1043 | 0.1077<br>0.1034 | 0.1075 | Ave | | 0.1070 | | | | 2.1 | | 15.0 | | | |
| Tert-amyl methyl ether | 0.6617<br>0.6962 | 0.6767<br>0.7331 | 0.7307<br>0.7077 | 0.7151<br>0.7140 | 0.6826 | Ave | | 0.7020 | | | | 3.5 | | 15.0 | | | |
| n-Butanol | 0.0138<br>0.0096 | 0.0095<br>0.0106 | 0.0097<br>0.0104 | 0.0114<br>0.0105 | 0.0089 | Ave | | 0.0105 | | | | 13.7 | | 15.0 | | | |
| Ethyl acrylate | 0.2316<br>0.2216 | 0.2219<br>0.2309 | 0.2178<br>0.2294 | 0.2151<br>0.2318 | 0.2122 | Ave | | 0.2236 | | | | 3.4 | | 15.0 | | | |
| Methyl methacrylate | 0.1524<br>0.1436 | 0.1363<br>0.1478 | 0.1383<br>0.1452 | 0.1404<br>0.1466 | 0.1384 | Ave | | 0.1432 | | | | 3.7 | | 15.0 | | | |
| 2-Nitropropane | +++++<br>0.0226 | 0.0209<br>0.0266 | 0.0198<br>0.0269 | 0.0195<br>0.0236 | 0.0194 | Ave | | 0.0224 | | | | 13.7 | | 15.0 | | | |
| n-Butyl acetate | +++++<br>0.3716 | +++++<br>0.3944 | 0.4011<br>0.3902 | 0.3921<br>0.3867 | 0.3596 | Ave | | 0.3851 | | | | 3.7 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50 Calibration End Date: 08/13/2015 20:42 Calibration ID: 30158

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1-Chlorohexane | 0.4929<br>0.4383 | 0.4320<br>0.4614 | 0.4729<br>0.4436 | 0.4373<br>0.4435 | 0.4315 | Ave | | 0.4504 | | | | 4.7 | | 15.0 | | | |
| Cyclohexanone | 0.0128<br>0.0108 | 0.0155<br>0.0119 | 0.0120<br>0.0114 | 0.0149<br>0.0114 | 0.0103 | Ave | | 0.0123 | | | | 14.4 | | 15.0 | | | |
| 1,2,3-Trimethylbenzene | 2.8189<br>3.0746 | 2.8992<br>3.2639 | 3.1500<br>3.1145 | 3.0040<br>3.0651 | 2.9168 | Ave | | 3.0341 | | | | 4.6 | | 15.0 | | | |
| 1,3,5-Trichlorobenzene | 1.1368<br>1.0908 | 1.0640<br>1.1383 | 1.1518<br>1.0486 | 1.1209<br>1.0391 | 1.0725 | Ave | | 1.0959 | | | | 3.9 | | 15.0 | | | |
| 2-Methylnaphthalene | 1.1258<br>1.2729 | 1.0980<br>1.4105 | 1.0984<br>1.2876 | 1.2519<br>1.2359 | 1.1966 | Ave | | 1.2197 | | | | 8.4 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50 Calibration End Date: 08/13/2015 20:42 Calibration ID: 30158

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-193150/20 | 144584.D |
| Level 2 | IC 240-193150/19 | 144583.D |
| Level 3 | IC 240-193150/18 | 144582.D |
| Level 4 | IC 240-193150/17 | 144581.D |
| Level 5 | IC 240-193150/16 | 144580.D |
| Level 6 | IC 240-193150/15 | 144579.D |
| Level 7 | IC 240-193150/14 | 144578.D |
| Level 8 | IC 240-193150/13 | 144577.D |
| Level 9 | IC 240-193150/12 | 144576.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Acetonitrile | FB | Ave | 7388<br>303462 | 10973<br>632599 | 30131<br>1256831 | 59051<br>1567842 | 126527 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| 2-Chloro-1,3-butadiene | FB | Ave | 10787<br>606721 | 22288<br>1265427 | 65473<br>2509599 | 121146<br>3096439 | 236791 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Diisopropyl ether | FB | Ave | 5444<br>292480 | 10602<br>607343 | 29596<br>1187824 | 60452<br>1459333 | 113175 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl t-butyl ether | FB | Ave | 22036<br>1141245 | 44869<br>2409991 | 114365<br>4840430 | 226764<br>5986416 | 436960 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl acetate | FB | Ave | 11385<br>435795 | 16786<br>895870 | 40612<br>1806750 | 84961<br>2228660 | 167431 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Propionitrile | FB | Ave | 7963<br>374531 | 16256<br>786705 | 37500<br>1563948 | 77984<br>1921197 | 145100 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| Methacrylonitrile | FB | Ave | 29273<br>1516021 | 58312<br>3095674 | 146830<br>5994798 | 301406<br>7302328 | 590265 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| Tert-amyl methyl ether | FB | Ave | 18089<br>975479 | 37373<br>2052126 | 98783<br>4066811 | 200083<br>5043426 | 374623 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| n-Butanol | CBZ | Ave | 5835<br>210549 | 8000<br>460443 | 20632<br>903775 | 49843<br>1135945 | 77408 | 25.0<br>1250 | 50.0<br>2500 | 125<br>5000 | 250<br>6250 | 500 |
| Ethyl acrylate | FB | Ave | 6332<br>310518 | 12255<br>646338 | 29451<br>1317957 | 60196<br>1637141 | 116465 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methyl methacrylate | FB | Ave | 8332<br>402313 | 15055<br>827639 | 37395<br>1668202 | 78576<br>2071659 | 151896 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| 2-Nitropropane | FB | Ave | +++++<br>63467 | 2313<br>149194 | 5344<br>309024 | 10920<br>333900 | 21282 | +++++<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| n-Butyl acetate | CBZ | Ave | +++++<br>325474 | +++++<br>683511 | 34073<br>1362203 | 68874<br>1673169 | 125276 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1-Chlorohexane | CBZ | Ave | 8362<br>383892 | 14627<br>799511 | 40171<br>1548771 | 76810<br>1919108 | 150333 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50 Calibration End Date: 08/13/2015 20:42 Calibration ID: 30158

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Cyclohexanone | CBZ | Ave | 2169<br>94475 | 5243<br>205876 | 10210<br>398420 | 26100<br>493296 | 35769 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| 1,2,3-Trimethylbenzene | DCB | Ave | 22714<br>1297881 | 46268<br>2748737 | 129836<br>5348625 | 256074<br>6584366 | 492514 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3,5-Trichlorobenzene | DCB | Ave | 9160<br>460463 | 16981<br>958626 | 47473<br>1800789 | 95549<br>2232228 | 181103 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Methylnaphthalene | DCB | Ave | 18143<br>1074702 | 35046<br>2375842 | 90547<br>4422384 | 213428<br>5309881 | 404111 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |

Curve Type Legend:
Ave = Average ISTD

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-196026/11 | 144943.D |
| Level 2 | IC 240-196026/10 | 144942.D |
| Level 3 | IC 240-196026/9 | 144941.D |
| Level 4 | IC 240-196026/8 | 144940.D |
| Level 5 | IC 240-196026/7 | 144939.D |
| Level 6 | ICIS 240-196026/6 | 144938.D |
| Level 7 | IC 240-196026/5 | 144937.D |
| Level 8 | IC 240-196026/4 | 144936.D |
| Level 9 | IC 240-196026/3 | 144935.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | +++++<br>0.2391 | +++++<br>0.2507 | 0.2470<br>0.2355 | 0.2306<br>0.2458 | 0.2468 | Ave | | 0.2422 | | | 0.1000 | 3.0 | | 15.0 | | | |
| Chloromethane | 0.3754<br>0.3823 | 0.4019<br>0.3850 | 0.3740<br>0.3707 | 0.3723<br>0.3714 | 0.3751 | Ave | | 0.3787 | | | 0.1000 | 2.6 | | 15.0 | | | |
| Vinyl chloride | 0.2535<br>0.2723 | 0.2872<br>0.2816 | 0.2779<br>0.2705 | 0.2715<br>0.2802 | 0.2773 | Ave | | 0.2747 | | | 0.1000 | 3.5 | | 15.0 | | | |
| Butadiene | 0.2421<br>0.3144 | 0.2898<br>0.3180 | 0.3006<br>0.3092 | 0.3016<br>0.3170 | 0.3102 | Ave | | 0.3003 | | | | 7.9 | | 15.0 | | | |
| Bromomethane | 0.1681<br>0.1234 | 0.1376<br>+++++ | 0.1308<br>+++++ | 0.1354<br>+++++ | 0.1348 | Ave | | 0.1383 | | | 0.0500 | 11.1 | | 15.0 | | | |
| Chloroethane | 0.1729<br>0.1392 | 0.1452<br>0.1270 | 0.1461<br>+++++ | 0.1422<br>+++++ | 0.1409 | Ave | | 0.1448 | | | | 9.6 | | 15.0 | | | |
| Dichlorofluoromethane | 0.3353<br>0.3214 | 0.3424<br>0.3238 | 0.3254<br>0.3013 | 0.3326<br>0.3083 | 0.3302 | Ave | | 0.3245 | | | | 4.0 | | 15.0 | | | |
| Trichlorofluoromethane | 0.2372<br>0.3085 | 0.2496<br>0.3130 | 0.2776<br>0.2931 | 0.2862<br>0.3038 | 0.2978 | Ave | | 0.2852 | | | 0.1000 | 9.2 | | 15.0 | | | |
| Ethyl ether | 0.2263<br>0.1633 | 0.1893<br>0.1738 | 0.1892<br>0.1620 | 0.1718<br>0.1640 | 0.1839 | Ave | | 0.1804 | | | | 11.2 | | 15.0 | | | |
| Acrolein | 0.0216<br>0.0171 | 0.0159<br>0.0175 | 0.0196<br>+++++ | 0.0166<br>+++++ | 0.0179 | Ave | | 0.0180 | | | | 10.9 | | 15.0 | | | |
| 1,1-Dichloroethene | 0.2079<br>0.2313 | 0.1781<br>0.2399 | 0.2268<br>0.2226 | 0.2292<br>0.2287 | 0.2305 | Ave | | 0.2217 | | | 0.1000 | 8.3 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.1337<br>0.1711 | 0.1398<br>0.1761 | 0.1720<br>0.1647 | 0.1739<br>0.1677 | 0.1631 | Ave | | 0.1625 | | | 0.0500 | 9.4 | | 15.0 | | | |
| Acetone | +++++<br>0.0712 | +++++<br>0.0711 | 0.1266<br>0.0622 | 0.0892<br>0.0631 | 0.0783 | Lin1 | 0.6353 | 0.0626 | | | 0.0500 | | | | 0.9980 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Iodomethane | 0.3031<br>0.3309 | 0.3161<br>0.3401 | 0.3156<br>0.3166 | 0.3107<br>0.3246 | 0.3168 | Ave | | 0.3194 | | | | 3.4 | | 15.0 | | | |
| Carbon disulfide | +++++<br>0.6938 | +++++<br>0.7470 | 0.5786<br>0.7111 | 0.6047<br>0.7368 | 0.6277 | Ave | | 0.6714 | | | 0.1000 | 10.0 | | 15.0 | | | |
| 3-Chloro-1-propene | +++++<br>0.1475 | +++++<br>0.1563 | 0.1275<br>0.1460 | 0.1372<br>0.1497 | 0.1368 | Ave | | 0.1430 | | | | 6.8 | | 15.0 | | | |
| Methyl acetate | 0.1615<br>0.1574 | 0.1594<br>0.1602 | 0.1513<br>0.1539 | 0.1584<br>0.1527 | 0.1554 | Ave | | 0.1567 | | | 0.1000 | 2.3 | | 15.0 | | | |
| Methylene Chloride | +++++<br>0.2588 | +++++<br>0.2535 | 0.5579<br>0.2285 | 0.3574<br>0.2329 | 0.3001 | Lin1 | 1.5559 | 0.2263 | | | 0.1000 | | | | 0.9990 | | 0.9900 |
| tert-Butyl alcohol (TBA) | 0.0223<br>0.0196 | 0.0192<br>0.0202 | 0.0186<br>0.0190 | 0.0196<br>0.0197 | 0.0203 | Ave | | 0.0198 | | | | 5.4 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.2358<br>0.2589 | 0.2329<br>0.2662 | 0.2511<br>0.2514 | 0.2599<br>0.2562 | 0.2510 | Ave | | 0.2515 | | | 0.1000 | 4.3 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.5407<br>0.5781 | 0.5287<br>0.5970 | 0.5676<br>0.5608 | 0.5729<br>0.5665 | 0.5729 | Ave | | 0.5650 | | | 0.1000 | 3.6 | | 15.0 | | | |
| Acrylonitrile | 0.0788<br>0.0761 | 0.0771<br>0.0770 | 0.0752<br>0.0728 | 0.0765<br>0.0724 | 0.0759 | Ave | | 0.0757 | | | | 2.7 | | 15.0 | | | |
| Hexane | +++++<br>0.0773 | +++++<br>0.0782 | 0.1040<br>0.0738 | 0.0811<br>0.0760 | 0.0806 | Ave | | 0.0816 | | | | 12.5 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.4672<br>0.5014 | 0.4509<br>0.5208 | 0.4998<br>0.4916 | 0.4999<br>0.5087 | 0.4920 | Ave | | 0.4925 | | | 0.1000 | 4.3 | | 15.0 | | | |
| Vinyl acetate | 0.0241<br>0.0308 | 0.0166<br>0.0307 | 0.0238<br>+++++ | 0.0268<br>+++++ | 0.0267 | Lin1 | -0.018 | 0.0303 | | | | | | | 0.9970 | | 0.9900 |
| 2,2-Dichloropropane | 0.1821<br>0.2315 | 0.1870<br>0.2370 | 0.2083<br>0.2070 | 0.2175<br>0.2121 | 0.2141 | Ave | | 0.2107 | | | | 8.5 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.2557<br>0.2665 | 0.2545<br>0.2811 | 0.2602<br>0.2612 | 0.2694<br>0.2684 | 0.2713 | Ave | | 0.2654 | | | 0.1000 | 3.2 | | 15.0 | | | |
| 2-Butanone (MEK) | +++++<br>0.0942 | 0.1118<br>0.0991 | 0.0952<br>0.0916 | 0.0894<br>0.0929 | 0.0959 | Ave | | 0.0962 | | | 0.0500 | 7.2 | | 15.0 | | | |
| Bromochloromethane | 0.0858<br>0.1003 | 0.0854<br>0.1062 | 0.0973<br>0.1000 | 0.0967<br>0.1004 | 0.0986 | Ave | | 0.0967 | | | | 7.1 | | 15.0 | | | |
| Tetrahydrofuran | +++++<br>0.0638 | 0.0727<br>0.0652 | 0.0687<br>0.0628 | 0.0716<br>0.0619 | 0.0652 | Ave | | 0.0665 | | | | 6.1 | | 15.0 | | | |
| Chloroform | 0.4109<br>0.3987 | 0.3844<br>0.4148 | 0.3916<br>0.3944 | 0.3857<br>0.4100 | 0.3904 | Ave | | 0.3979 | | | | 2.9 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.2815<br>0.3335 | 0.2455<br>0.3520 | 0.3014<br>0.3288 | 0.3164<br>0.3392 | 0.3164 | Ave | | 0.3127 | | | 0.1000 | 10.5 | | 15.0 | | | |
| Cyclohexane | 0.5387<br>0.6083 | 0.4961<br>0.6348 | 0.6113<br>0.5854 | 0.6257<br>0.6043 | 0.6056 | Ave | | 0.5900 | | | 0.1000 | 7.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon tetrachloride | 0.1974<br>0.2765 | 0.2070<br>0.2949 | 0.2344<br>0.2806 | 0.2508<br>0.2887 | 0.2562 | Ave | | 0.2541 | | | 0.1000 | 13.8 | | 15.0 | | | |
| 1,1-Dichloropropene | +++++<br>0.3352 | +++++<br>0.3502 | 0.3354<br>0.3258 | 0.3276<br>0.3378 | 0.3200 | Ave | | 0.3331 | | | | 3.0 | | 15.0 | | | |
| Benzene | 0.9461<br>1.0076 | 0.9369<br>1.0637 | 0.9831<br>0.9992 | 1.0297<br>1.0364 | 0.9873 | Ave | | 0.9989 | | | 0.5000 | 4.1 | | 15.0 | | | |
| Isobutyl alcohol | 0.0079<br>0.0070 | 0.0073<br>0.0070 | 0.0070<br>0.0065 | 0.0074<br>0.0064 | 0.0078 | Ave | | 0.0071 | | | | 7.3 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.2922<br>0.2806 | 0.2538<br>0.2902 | 0.2809<br>0.2746 | 0.2857<br>0.2775 | 0.2775 | Ave | | 0.2792 | | | 0.1000 | 4.0 | | 15.0 | | | |
| n-Heptane | 0.0495<br>0.0719 | 0.0549<br>0.0771 | 0.0671<br>0.0732 | 0.0733<br>0.0762 | 0.0704 | Ave | | 0.0682 | | | | 14.1 | | 15.0 | | | |
| Trichloroethene | 0.2322<br>0.2352 | 0.1925<br>0.2515 | 0.2312<br>0.2339 | 0.2474<br>0.2400 | 0.2383 | Ave | | 0.2336 | | | 0.2000 | 7.2 | | 15.0 | | | |
| Methylcyclohexane | 0.3977<br>0.5007 | 0.4180<br>0.5258 | 0.5131<br>0.4892 | 0.4996<br>0.5023 | 0.4970 | Ave | | 0.4826 | | | 0.1000 | 9.1 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.2351<br>0.2464 | 0.2304<br>0.2703 | 0.2458<br>0.2552 | 0.2519<br>0.2638 | 0.2445 | Ave | | 0.2493 | | | 0.1000 | 5.1 | | 15.0 | | | |
| Dibromomethane | 0.1138<br>0.1049 | 0.0928<br>0.1099 | 0.1030<br>0.1045 | 0.0989<br>0.1052 | 0.1030 | Ave | | 0.1040 | | | | 5.7 | | 15.0 | | | |
| 1,4-Dioxane | 0.0016<br>0.0015 | 0.0018<br>0.0016 | 0.0017<br>0.0016 | 0.0016<br>0.0016 | 0.0018 | Ave | | 0.0017 | | | | 5.4 | | 15.0 | | | |
| Bromodichloromethane | +++++<br>0.2583 | +++++<br>0.2789 | 0.2043<br>0.2740 | 0.2214<br>0.2839 | 0.2350 | Ave | | 0.2508 | | | 0.1000 | 12.3 | | 15.0 | | | |
| 2-Chloroethyl vinyl ether | 0.0880<br>0.0981 | 0.0738<br>0.1047 | 0.0849<br>0.1021 | 0.0908<br>0.1040 | 0.0933 | Ave | | 0.0933 | | | | 10.9 | | 15.0 | | | |
| cis-1,3-Dichloropropene | +++++<br>0.3137 | +++++<br>0.3452 | 0.2529<br>0.3385 | 0.2764<br>0.3487 | 0.2874 | Ave | | 0.3090 | | | 0.1500 | 12.2 | | 15.0 | | | |
| 4-Methyl-2-pentanone (MIBK) | 0.3475<br>0.3097 | 0.3025<br>0.3146 | 0.3102<br>0.2962 | 0.3082<br>0.2899 | 0.3133 | Ave | | 0.3102 | | | 0.0500 | 5.2 | | 15.0 | | | |
| Toluene | 1.6889<br>1.6108 | 1.5730<br>1.6794 | 1.5789<br>1.5858 | 1.6326<br>1.6292 | 1.5722 | Ave | | 1.6167 | | | 0.4000 | 2.8 | | 15.0 | | | |
| trans-1,3-Dichloropropene | +++++<br>0.3786 | +++++<br>0.4056 | 0.2981<br>0.3958 | 0.3438<br>0.4021 | 0.3521 | Ave | | 0.3680 | | | 0.1000 | 10.6 | | 15.0 | | | |
| Ethyl methacrylate | 0.3411<br>0.2949 | 0.2728<br>0.3134 | 0.2921<br>0.2986 | 0.3041<br>0.3000 | 0.2937 | Ave | | 0.3012 | | | | 6.1 | | 15.0 | | | |
| 1,1,2-Trichloroethane | 0.2387<br>0.2409 | 0.2020<br>0.2502 | 0.2341<br>0.2296 | 0.2342<br>0.2336 | 0.2367 | Ave | | 0.2333 | | | 0.1000 | 5.6 | | 15.0 | | | |
| Tetrachloroethene | 0.2604<br>0.2924 | 0.2503<br>0.3024 | 0.2921<br>0.2817 | 0.2907<br>0.2908 | 0.2885 | Ave | | 0.2833 | | | 0.2000 | 6.0 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,3-Dichloropropane | 0.4985<br>0.4202 | 0.3942<br>0.4386 | 0.3932<br>0.4193 | 0.4331<br>0.4249 | 0.4155 | Ave | | 0.4264 | | | | 7.3 | | 15.0 | | | |
| 2-Hexanone | 0.2259<br>0.1948 | 0.2056<br>0.1966 | 0.2059<br>0.1806 | 0.2139<br>0.1792 | 0.2041 | Ave | | 0.2007 | | | 0.0500 | 7.5 | | 15.0 | | | |
| Dibromochloromethane | +++++<br>0.2379 | 0.1676<br>0.2584 | 0.1731<br>0.2529 | 0.1970<br>0.2554 | 0.2047 | Lin1 | -0.324 | 0.2543 | | | | | | | 0.9990 | | 0.9900 |
| 1,2-Dibromoethane | 0.1915<br>0.2254 | 0.1731<br>0.2352 | 0.2140<br>0.2225 | 0.2199<br>0.2231 | 0.2195 | Ave | | 0.2138 | | | | 9.0 | | 15.0 | | | |
| Chlorobenzene | 0.8957<br>0.9330 | 0.9144<br>0.9681 | 0.9220<br>0.9247 | 0.9499<br>0.9450 | 0.9159 | Ave | | 0.9298 | | | 0.3000 | 2.3 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.2424<br>0.3183 | 0.2461<br>0.3292 | 0.2688<br>0.3070 | 0.2915<br>0.3071 | 0.3010 | Ave | | 0.2902 | | | | 10.7 | | 15.0 | | | |
| Ethylbenzene | 0.4967<br>0.5429 | 0.5117<br>0.5557 | 0.5343<br>0.5186 | 0.5496<br>0.5291 | 0.5275 | Ave | | 0.5296 | | | | 3.5 | | 15.0 | | | |
| m-Xylene & p-Xylene | 0.6493<br>0.6610 | 0.5935<br>0.6804 | 0.6630<br>0.6340 | 0.6529<br>0.6481 | 0.6392 | Ave | | 0.6468 | | | | 3.7 | | 15.0 | | | |
| o-Xylene | 0.6484<br>0.6613 | 0.5588<br>0.6783 | 0.6201<br>0.6144 | 0.6581<br>0.6145 | 0.6575 | Ave | | 0.6346 | | | | 5.7 | | 15.0 | | | |
| Styrene | 0.9630<br>0.9968 | 0.8800<br>1.0308 | 0.8828<br>0.9707 | 0.9401<br>0.9819 | 0.9540 | Ave | | 0.9556 | | | 0.3000 | 5.2 | | 15.0 | | | |
| Bromoform | 0.0617<br>0.1189 | 0.0736<br>0.1321 | 0.0749<br>0.1315 | 0.0937<br>0.1318 | 0.0977 | Lin1 | -0.136 | 0.1302 | | | 0.1000 | | | | 0.9970 | | 0.9900 |
| Isopropylbenzene | 1.6615<br>1.8186 | 1.5426<br>1.8815 | 1.7312<br>1.7266 | 1.7956<br>1.7460 | 1.7709 | Ave | | 1.7416 | | | 0.1000 | 5.6 | | 15.0 | | | |
| Bromobenzene | 0.7018<br>0.7269 | 0.6199<br>0.7522 | 0.6741<br>0.7263 | 0.7025<br>0.7372 | 0.6965 | Ave | | 0.7042 | | | | 5.6 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 0.5544<br>0.6344 | 0.5565<br>0.6397 | 0.5491<br>0.6163 | 0.5930<br>0.5999 | 0.6184 | Ave | | 0.5958 | | | 0.3000 | 5.9 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.1905<br>0.1847 | 0.1628<br>0.1813 | 0.1632<br>0.1748 | 0.1767<br>0.1723 | 0.1805 | Ave | | 0.1763 | | | | 5.3 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | 0.2675<br>0.2192 | 0.1583<br>0.2222 | 0.1934<br>0.2263 | 0.2038<br>0.2273 | 0.2186 | Ave | | 0.2152 | | | | 13.7 | | 15.0 | | | |
| N-Propylbenzene | 0.9746<br>1.0339 | 0.8541<br>1.0565 | 0.8793<br>1.0140 | 0.9898<br>1.0407 | 0.9489 | Ave | | 0.9769 | | | | 7.3 | | 15.0 | | | |
| 2-Chlorotoluene | 0.8414<br>0.8319 | 0.6742<br>0.8400 | 0.7804<br>0.8160 | 0.8063<br>0.8183 | 0.8038 | Ave | | 0.8014 | | | | 6.4 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 2.7121<br>3.3648 | 2.7168<br>3.3834 | 2.9008<br>3.2276 | 3.1601<br>3.2885 | 3.1601 | Ave | | 3.1016 | | | | 8.4 | | 15.0 | | | |
| 4-Chlorotoluene | 0.8033<br>0.7833 | 0.7373<br>0.8244 | 0.7337<br>0.7938 | 0.7735<br>0.8011 | 0.7475 | Ave | | 0.7776 | | | | 4.1 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| tert-Butylbenzene | 2.3374<br>2.9631 | 2.4133<br>3.0556 | 2.6550<br>2.8376 | 2.8314<br>2.8634 | 2.8090 | Ave | | 2.7518 | | | | 8.7 | | 15.0 | | | |
| 1,2,4-Trimethylbenzene | 2.7336<br>3.2868 | 2.6205<br>3.3941 | 2.8875<br>3.2352 | 3.0975<br>3.2908 | 3.1820 | Ave | | 3.0809 | | | | 8.8 | | 15.0 | | | |
| sec-Butylbenzene | 3.5842<br>4.5688 | 3.5902<br>4.6258 | 3.9612<br>4.3978 | 4.2686<br>4.4271 | 4.3250 | Ave | | 4.1943 | | | | 9.4 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.4872<br>1.4944 | 1.4279<br>1.5057 | 1.4500<br>1.4639 | 1.4918<br>1.4763 | 1.4693 | Ave | | 1.4741 | | | 0.6000 | 1.7 | | 15.0 | | | |
| p-Isopropyltoluene | 2.6700<br>3.5326 | 2.7312<br>3.5868 | 3.1681<br>3.3752 | 3.3032<br>3.4312 | 3.3784 | Ave | | 3.2418 | | | | 10.2 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 1.6796<br>1.4771 | 1.4525<br>1.4817 | 1.3997<br>1.4447 | 1.4789<br>1.4779 | 1.4661 | Ave | | 1.4842 | | | 0.5000 | 5.2 | | 15.0 | | | |
| n-Butylbenzene | 2.6703<br>3.3398 | 2.7947<br>3.4180 | 3.0393<br>3.3135 | 3.1433<br>3.3757 | 3.2352 | Ave | | 3.1478 | | | | 8.4 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.5577<br>1.4312 | 1.4056<br>1.4532 | 1.3574<br>1.3836 | 1.4146<br>1.3837 | 1.4022 | Ave | | 1.4210 | | | 0.4000 | 4.1 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | 0.0691<br>0.0964 | 0.0727<br>0.1007 | 0.0775<br>0.1005 | 0.0817<br>0.0987 | 0.0866 | Ave | | 0.0871 | | | 0.0500 | 14.3 | | 15.0 | | | |
| 1,2,4-Trichlorobenzene | 0.9903<br>0.9952 | 0.9066<br>0.9801 | 0.9495<br>0.9392 | 1.0303<br>0.9542 | 1.0144 | Ave | | 0.9733 | | | 0.2000 | 4.0 | | 15.0 | | | |
| Hexachlorobutadiene | 0.3960<br>0.5853 | 0.4364<br>0.5796 | 0.5790<br>0.5455 | 0.5879<br>0.5625 | 0.5837 | Ave | | 0.5396 | | | | 13.3 | | 15.0 | | | |
| Naphthalene | 2.5487<br>2.3198 | 2.2039<br>2.2590 | 2.1592<br>2.1437 | 2.2755<br>2.0812 | 2.3749 | Ave | | 2.2629 | | | | 6.2 | | 15.0 | | | |
| 1,2,3-Trichlorobenzene | 0.9121<br>0.9513 | 0.9280<br>0.9117 | 0.9440<br>0.8603 | 0.9768<br>0.8520 | 0.9646 | Ave | | 0.9223 | | | | 4.7 | | 15.0 | | | |
| Dibromofluoromethane (Surr) | +++++<br>0.2047 | +++++<br>0.2121 | 0.2454<br>0.1982 | 0.2289<br>0.2019 | 0.2161 | Ave | | 0.2154 | | | | 7.8 | | 15.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | +++++<br>0.2276 | +++++<br>0.2392 | 0.3027<br>0.2282 | 0.2781<br>0.2330 | 0.2521 | Ave | | 0.2516 | | | | 11.4 | | 15.0 | | | |
| Toluene-d8 (Surr) | +++++<br>1.4443 | +++++<br>1.4774 | 1.8704<br>1.3884 | 1.6510<br>1.4120 | 1.5274 | Ave | | 1.5387 | | | | 11.1 | | 15.0 | | | |
| 4-Bromofluorobenzene (Surr) | +++++<br>1.0832 | +++++<br>1.0727 | 1.4341<br>1.0549 | 1.2044<br>1.0662 | 1.1169 | Ave | | 1.1475 | | | | 11.9 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-196026/11 | 144943.D |
| Level 2 | IC 240-196026/10 | 144942.D |
| Level 3 | IC 240-196026/9 | 144941.D |
| Level 4 | IC 240-196026/8 | 144940.D |
| Level 5 | IC 240-196026/7 | 144939.D |
| Level 6 | ICIS 240-196026/6 | 144938.D |
| Level 7 | IC 240-196026/5 | 144937.D |
| Level 8 | IC 240-196026/4 | 144936.D |
| Level 9 | IC 240-196026/3 | 144935.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Dichlorodifluoromethane | FB | Ave | +++++<br>356137 | +++++<br>748933 | 36569<br>1466377 | 68169<br>1835598 | 147021 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Chloromethane | FB | Ave | 10383<br>569551 | 22625<br>1150096 | 55373<br>2307997 | 110043<br>2773342 | 223410 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Vinyl chloride | FB | Ave | 7011<br>405713 | 16172<br>841040 | 41149<br>1684326 | 80229<br>2092495 | 165146 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Butadiene | FB | Ave | 6696<br>468393 | 16314<br>949979 | 44512<br>1925445 | 89146<br>2367224 | 184770 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Bromomethane | FB | Ave | 4648<br>183791 | 7747<br>+++++ | 19368<br>+++++ | 40006<br>+++++ | 80268 | 1.00<br>50.0 | 2.00<br>+++++ | 5.00<br>+++++ | 10.0<br>+++++ | 20.0 |
| Chloroethane | FB | Ave | 4781<br>207308 | 8175<br>379413 | 21638<br>+++++ | 42019<br>+++++ | 83922 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>+++++ | 10.0<br>+++++ | 20.0 |
| Dichlorofluoromethane | FB | Ave | 9272<br>478742 | 19276<br>967265 | 48180<br>1875722 | 98295<br>2301848 | 196686 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Trichlorofluoromethane | FB | Ave | 6559<br>459614 | 14053<br>934958 | 41107<br>1824715 | 84588<br>2268583 | 177390 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl ether | FB | Ave | 6258<br>243336 | 10659<br>519027 | 28019<br>1008376 | 50769<br>1224410 | 109526 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Acrolein | FB | Ave | 2993<br>127696 | 4483<br>261405 | 14479<br>+++++ | 24467<br>+++++ | 53257 | 5.00<br>250 | 10.0<br>500 | 25.0<br>+++++ | 50.0<br>+++++ | 100 |
| 1,1-Dichloroethene | FB | Ave | 5750<br>344590 | 10027<br>716734 | 33584<br>1386108 | 67737<br>1707262 | 137258 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroetha ne | FB | Ave | 3699<br>254932 | 7868<br>526084 | 25468<br>1025667 | 51393<br>1251857 | 97164 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Acetone | FB | Lin1 | +++++<br>212159 | +++++<br>424676 | 37501<br>774275 | 52703<br>941871 | 93287 | +++++<br>100 | +++++<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Iodomethane | FB | Ave | 8383<br>493006 | 17794<br>1015933 | 46725<br>1971089 | 91832<br>2423809 | 188704 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Carbon disulfide | FB | Ave | +++++<br>1033515 | +++++<br>2231244 | 85670<br>4427381 | 178714<br>5501396 | 373862 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 3-Chloro-1-propene | FB | Ave | +++++<br>219767 | +++++<br>467005 | 18885<br>908999 | 40539<br>1117889 | 81474 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methyl acetate | FB | Ave | 22330<br>1172266 | 44885<br>2392801 | 112036<br>4789789 | 234059<br>5701867 | 462647 | 5.00<br>250 | 10.0<br>500 | 25.0<br>1000 | 50.0<br>1250 | 100 |
| Methylene Chloride | FB | Lin1 | +++++<br>385484 | +++++<br>757088 | 82611<br>1422440 | 105644<br>1738961 | 178739 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| tert-Butyl alcohol (TBA) | FB | Ave | 6178<br>292151 | 10809<br>601929 | 27588<br>1184581 | 57843<br>1471581 | 120716 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| trans-1,2-Dichloroethene | FB | Ave | 6522<br>385670 | 13114<br>795293 | 37178<br>1565230 | 76822<br>1912734 | 149499 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methyl tert-butyl ether | FB | Ave | 14953<br>861134 | 29764<br>1783409 | 84049<br>3491615 | 169317<br>4230179 | 341246 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Acrylonitrile | FB | Ave | 21785<br>1133959 | 43399<br>2301252 | 111279<br>4534284 | 225966<br>5402864 | 451855 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| Hexane | FB | Ave | +++++<br>115126 | +++++<br>233508 | 15392<br>459224 | 23963<br>567539 | 48009 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1-Dichloroethane | FB | Ave | 12922<br>746990 | 25387<br>1555732 | 74007<br>3060770 | 147740<br>3798439 | 293058 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Vinyl acetate | FB | Lin1 | 667<br>45823 | 932<br>91562 | 3519<br>+++++ | 7932<br>+++++ | 15884 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>+++++ | 10.0<br>+++++ | 20.0 |
| 2,2-Dichloropropane | FB | Ave | 5036<br>344804 | 10529<br>708031 | 30835<br>1289042 | 64289<br>1583856 | 127532 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| cis-1,2-Dichloroethene | FB | Ave | 7073<br>396968 | 14329<br>839540 | 38524<br>1626156 | 79621<br>2004348 | 161592 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Butanone (MEK) | FB | Ave | +++++<br>280552 | 12585<br>591753 | 28186<br>1140829 | 52850<br>1386909 | 114195 | +++++<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Bromochloromethane | FB | Ave | 2372<br>149423 | 4806<br>317360 | 14400<br>622349 | 28593<br>749740 | 58707 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Tetrahydrofuran | FB | Ave | +++++<br>189961 | 8190<br>389802 | 20359<br>781838 | 42328<br>924687 | 77626 | +++++<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Chloroform | FB | Ave | 11364<br>593941 | 21643<br>1239162 | 57984<br>2455954 | 113989<br>3061654 | 232521 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,1-Trichloroethane | FB | Ave | 7784<br>496863 | 13820<br>1051452 | 44626<br>2047378 | 93504<br>2532683 | 188459 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Cyclohexane | FB | Ave | 14900<br>906203 | 27932<br>1896235 | 90514<br>3644821 | 184933<br>4512267 | 360725 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Carbon tetrachloride | FB | Ave | 5460<br>411902 | 11655<br>880737 | 34704<br>1747228 | 74120<br>2155848 | 152621 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1-Dichloropropene | FB | Ave | +++++<br>499401 | +++++<br>1046178 | 49664<br>2028361 | 96828<br>2521880 | 190594 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Benzene | FB | Ave | 26167<br>1500996 | 52745<br>3177421 | 145567<br>6221614 | 304344<br>7738061 | 588026 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Isobutyl alcohol | FB | Ave | 5460<br>261074 | 10213<br>525940 | 25933<br>1010906 | 54805<br>1189785 | 116205 | 25.0<br>1250 | 50.0<br>2500 | 125<br>5000 | 250<br>6250 | 500 |
| 1,2-Dichloroethane | FB | Ave | 8081<br>418078 | 14287<br>866965 | 41595<br>1709845 | 84448<br>2072165 | 165251 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| n-Heptane | FB | Ave | 1370<br>107045 | 3089<br>230164 | 9941<br>455483 | 21652<br>568917 | 41942 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Trichloroethene | FB | Ave | 6423<br>350358 | 10835<br>751379 | 34227<br>1456606 | 73115<br>1792315 | 141942 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methylcyclohexane | FB | Ave | 10999<br>745915 | 23534<br>1570602 | 75969<br>3045944 | 147660<br>3750421 | 296027 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dichloropropane | FB | Ave | 6501<br>367015 | 12972<br>807304 | 36398<br>1589013 | 74445<br>1969959 | 145632 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Dibromomethane | FB | Ave | 3146<br>156268 | 5223<br>328160 | 15254<br>650540 | 29240<br>785166 | 61356 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,4-Dioxane | FB | Ave | 896<br>45240 | 2050<br>96254 | 4894<br>200368 | 9657<br>245495 | 21042 | 20.0<br>1000 | 40.0<br>2000 | 100<br>4000 | 200<br>5000 | 400 |
| Bromodichloromethane | FB | Ave | +++++<br>384799 | +++++<br>832998 | 30251<br>1706175 | 65422<br>2119797 | 139991 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Chloroethyl vinyl ether | FB | Ave | 4867<br>292355 | 8305<br>625520 | 25131<br>1271843 | 53690<br>1553516 | 111110 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| cis-1,3-Dichloropropene | FB | Ave | +++++<br>467262 | +++++<br>1031197 | 37447<br>2107636 | 81678<br>2603965 | 171161 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | CBZ | Ave | 11429<br>563617 | 20282<br>1178361 | 55788<br>2332635 | 109762<br>2766526 | 227626 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Toluene | CBZ | Ave | 27774<br>1465877 | 52739<br>3144782 | 141953<br>6243364 | 290701<br>7772646 | 571155 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| trans-1,3-Dichloropropene | CBZ | Ave | +++++<br>344567 | +++++<br>759613 | 26806<br>1558216 | 61223<br>1918393 | 127919 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl methacrylate | CBZ | Ave | 5609<br>268334 | 9146<br>586878 | 26258<br>1175568 | 54145<br>1431045 | 106690 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,2-Trichloroethane | CBZ | Ave | 3926<br>219261 | 6772<br>468456 | 21044<br>904034 | 41703<br>1114274 | 86000 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Tetrachloroethene | CBZ | Ave | 4283<br>266087 | 8391<br>566338 | 26258<br>1109089 | 51758<br>1387568 | 104805 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3-Dichloropropane | CBZ | Ave | 8198<br>382372 | 13218<br>821259 | 35354<br>1650749 | 77111<br>2027259 | 150930 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Hexanone | CBZ | Ave | 7429<br>354469 | 13785<br>736413 | 37025<br>1421673 | 76163<br>1709783 | 148306 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Dibromochloromethane | CBZ | Lin1 | +++++<br>216458 | 5618<br>483817 | 15565<br>995515 | 35077<br>1218638 | 74370 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1,2-Dibromoethane | CBZ | Ave | 3150<br>205128 | 5804<br>440477 | 19238<br>875919 | 39164<br>1064183 | 79732 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Chlorobenzene | CBZ | Ave | 14730<br>849090 | 30658<br>1812785 | 82892<br>3640404 | 169143<br>4508577 | 332714 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,1,2-Tetrachloroethane | CBZ | Ave | 3986<br>289664 | 8252<br>616483 | 24169<br>1208539 | 51910<br>1465339 | 109363 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethylbenzene | CBZ | Ave | 8169<br>494070 | 17157<br>1040646 | 48037<br>2041599 | 97870<br>2524079 | 191624 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| m-Xylene & p-Xylene | CBZ | Ave | 10677<br>601578 | 19898<br>1274050 | 59612<br>2495890 | 116259<br>3091955 | 232208 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| o-Xylene | CBZ | Ave | 10663<br>601783 | 18737<br>1270093 | 55753<br>2418978 | 117189<br>2931823 | 238863 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Styrene | CBZ | Ave | 15837<br>907146 | 29506<br>1930215 | 79372<br>3821764 | 167399<br>4684740 | 346563 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Bromoform | CBZ | Lin1 | 1015<br>108179 | 2467<br>247433 | 6730<br>517661 | 16693<br>628700 | 35497 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Isopropylbenzene | CBZ | Ave | 27324<br>1654997 | 51721<br>3523233 | 155645<br>6797585 | 319732<br>8330012 | 643348 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Bromobenzene | DCB | Ave | 5573<br>295347 | 9956<br>626961 | 28765<br>1217640 | 58476<br>1492683 | 116607 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,2,2-Tetrachloroethane | DCB | Ave | 4403<br>257748 | 8937<br>533250 | 23432<br>1033232 | 49363<br>1214765 | 103530 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,3-Trichloropropane | DCB | Ave | 1513<br>75031 | 2614<br>151158 | 6965<br>293104 | 14705<br>348980 | 30218 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| trans-1,4-Dichloro-2-butene | DCB | Ave | 2124<br>89070 | 2543<br>185204 | 8252<br>379325 | 16966<br>460351 | 36604 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| N-Propylbenzene | DCB | Ave | 7740<br>420052 | 13717<br>880640 | 37524<br>1699968 | 82391<br>2107190 | 158855 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Chlorotoluene | DCB | Ave | 6682<br>337984 | 10827<br>700147 | 33303<br>1368070 | 67113<br>1656963 | 134563 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3,5-Trimethylbenzene | DCB | Ave | 21538<br>1367130 | 43631<br>2820139 | 123786<br>5410976 | 263032<br>6658839 | 529049 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 4-Chlorotoluene | DCB | Ave | 6379<br>318271 | 11841<br>687196 | 31307<br>1330849 | 64384<br>1622168 | 125148 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| tert-Butylbenzene | DCB | Ave | 18562<br>1203890 | 38757<br>2546917 | 113296<br>4757117 | 235674<br>5798117 | 470266 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,4-Trimethylbenzene | DCB | Ave | 21708<br>1335433 | 42084<br>2829044 | 123218<br>5423776 | 257826<br>6663540 | 532704 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| sec-Butylbenzene | DCB | Ave | 28463<br>1856287 | 57658<br>3855727 | 169034<br>7372707 | 355305<br>8964364 | 724063 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3-Dichlorobenzene | DCB | Ave | 11810<br>607174 | 22932<br>1255024 | 61875<br>2454144 | 124176<br>2989387 | 245980 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| p-Isopropyltoluene | DCB | Ave | 21203<br>1435272 | 43863<br>2989705 | 135190<br>5658412 | 274948<br>6947777 | 565586 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,4-Dichlorobenzene | DCB | Ave | 13338<br>600156 | 23326<br>1235039 | 59729<br>2421962 | 123097<br>2992653 | 245441 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| n-Butylbenzene | DCB | Ave | 21206<br>1356939 | 44882<br>2849004 | 129694<br>5555001 | 261638<br>6835319 | 541621 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dichlorobenzene | DCB | Ave | 12370<br>581486 | 22574<br>1211298 | 57923<br>2319608 | 117745<br>2801754 | 234745 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | DCB | Ave | 549<br>39179 | 1167<br>83939 | 3309<br>168459 | 6800<br>199946 | 14490 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,4-Trichlorobenzene | DCB | Ave | 7864<br>404356 | 14560<br>816947 | 40518<br>1574505 | 85761<br>1932210 | 169826 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Hexachlorobutadiene | DCB | Ave | 3145<br>237801 | 7009<br>483093 | 24708<br>914525 | 48937<br>1138913 | 97726 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Naphthalene | DCB | Ave | 20240<br>942540 | 35394<br>1882955 | 92141<br>3593926 | 189406<br>4214177 | 397593 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,3-Trichlorobenzene | DCB | Ave | 7243<br>386525 | 14903<br>759901 | 40282<br>1442298 | 81306<br>1725103 | 161494 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Dibromofluoromethane (Surr) | FB | Ave | +++++<br>304994 | +++++<br>633562 | 36336<br>1234341 | 67662<br>1507714 | 128718 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Ave | +++++<br>339100 | +++++<br>714630 | 44817<br>1421037 | 82196<br>1739346 | 150166 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Toluene-d8 (Surr) | CBZ | Ave | +++++<br>1314337 | +++++<br>2766497 | 168163<br>5466311 | 293978<br>6736325 | 554888 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 4-Bromofluorobenzene (Surr) | DCB | Ave | +++++<br>440117 | +++++<br>894108 | 61195<br>1768529 | 100246<br>2158955 | 186984 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin1 = Linear 1/conc ISTD |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 189147

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/15/2015 01:51 Calibration End Date: 07/15/2015 03:38 Calibration ID: 29499

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STDA9 240-189147/9 | UX88550.D |
| Level 2 | STDA9 240-189147/8 | UX88549.D |
| Level 3 | STDA9 240-189147/7 | UX88548.D |
| Level 4 | STDA9 240-189147/6 | UX88547.D |
| Level 5 | STDA9 240-189147/5 | UX88546.D |
| Level 6 | STDA9 240-189147/4 | UX88545.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Acetonitrile | 0.0608 0.0519 | 0.0530 | 0.0453 | 0.0535 | 0.0494 | Ave | | 0.0523 | | | | 9.8 | | 15.0 | | | |
| Isopropyl ether | 0.2111 0.2108 | 0.2141 | 0.1940 | 0.2153 | 0.2219 | Ave | | 0.2112 | | | | 4.4 | | 15.0 | | | |
| 2-Chloro-1,3-butadiene | 0.5517 0.5693 | 0.5194 | 0.4971 | 0.5227 | 0.5605 | Ave | | 0.5368 | | | | 5.2 | | 15.0 | | | |
| Tert-butyl ethyl ether | 0.5272 0.6276 | 0.5357 | 0.5209 | 0.6108 | 0.6398 | Ave | | 0.5770 | | | | 9.5 | | 15.0 | | | |
| Ethyl acetate | 0.3474 0.3615 | 0.3251 | 0.3129 | 0.3675 | 0.3570 | Ave | | 0.3452 | | | | 6.3 | | 15.0 | | | |
| Propionitrile | 0.0595 0.0533 | 0.0527 | 0.0465 | 0.0552 | 0.0515 | Ave | | 0.0531 | | | | 8.1 | | 15.0 | | | |
| Methacrylonitrile | 0.2978 0.2716 | 0.2652 | 0.2431 | 0.2765 | 0.2694 | Ave | | 0.2706 | | | | 6.5 | | 15.0 | | | |
| Tert-amyl methyl ether | 0.4190 0.4932 | 0.3977 | 0.3839 | 0.4571 | 0.4777 | Ave | | 0.4381 | | | | 10.2 | | 15.0 | | | |
| n-Butanol | 0.0085 0.0084 | 0.0063 | 0.0049 | 0.0076 | 0.0072 | Qua | -0.067 | 0.0065 | 0.0000004 | | | | | | 0.9990 | | 0.9900 |
| Ethyl acrylate | 0.3465 0.4122 | 0.3263 | 0.3464 | 0.4078 | 0.4207 | Ave | | 0.3766 | | | | 11.0 | | 15.0 | | | |
| Methyl methacrylate | 0.3428 0.3586 | 0.3255 | 0.3165 | 0.3614 | 0.3716 | Ave | | 0.3461 | | | | 6.3 | | 15.0 | | | |
| 2-Nitropropane | 0.0560 0.0992 | 0.0577 | 0.0545 | 0.0829 | 0.0884 | Qua | -0.574 | 0.0818 | 0.0000471 | | | | | | 1.0000 | | 0.9900 |
| n-Butyl acetate | 0.5759 0.5133 | 0.4983 | 0.4371 | 0.5379 | 0.5473 | Ave | | 0.5183 | | | | 9.3 | | 15.0 | | | |
| 1-Chlorohexane | 0.4000 0.4054 | 0.3781 | 0.3626 | 0.3734 | 0.4152 | Ave | | 0.3891 | | | | 5.3 | | 15.0 | | | |
| Cyclohexanone | 0.0230 0.0165 | 0.0173 | 0.0121 | 0.0165 | 0.0149 | Qua | 0.3064 | 0.0136 | 0.0000013 | | | | | | 0.9980 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 189147

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/15/2015 01:51 Calibration End Date: 07/15/2015 03:38 Calibration ID: 29499

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,2,3-Trimethylbenzene | 2.9402 2.9004 | 2.7999 | 2.6150 | 2.7114 | 2.9307 | Ave | | 2.8163 | | | | 4.7 | | 15.0 | | | |
| 1,3,5-Trichlorobenzene | 1.2299 1.1578 | 1.1546 | 1.0879 | 1.0955 | 1.1550 | Ave | | 1.1468 | | | | 4.5 | | 15.0 | | | |
| 2-Methylnaphthalene | 2.0008 1.3028 | 1.6904 | 1.5655 | 1.5876 | 1.5761 | Ave | | 1.6206 | | | | 14.0 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 189147

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/15/2015 01:51 Calibration End Date: 07/15/2015 03:38 Calibration ID: 29499

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STDA9 240-189147/9 | UX88550.D |
| Level 2 | STDA9 240-189147/8 | UX88549.D |
| Level 3 | STDA9 240-189147/7 | UX88548.D |
| Level 4 | STDA9 240-189147/6 | UX88547.D |
| Level 5 | STDA9 240-189147/5 | UX88546.D |
| Level 6 | STDA9 240-189147/4 | UX88545.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Acetonitrile | FB | Ave | 54039<br>1946452 | 95176 | 170154 | 485511 | 904712 | 50.0<br>2000 | 100 | 200 | 500 | 1000 |
| Isopropyl ether | FB | Ave | 18760<br>791322 | 38415 | 72928 | 195279 | 406204 | 5.00<br>200 | 10.0 | 20.0 | 50.0 | 100 |
| 2-Chloro-1,3-butadiene | FB | Ave | 49023<br>2136848 | 93194 | 186816 | 474070 | 1026070 | 5.00<br>200 | 10.0 | 20.0 | 50.0 | 100 |
| Tert-butyl ethyl ether | FB | Ave | 46844<br>2355578 | 96118 | 195770 | 553989 | 1171169 | 5.00<br>200 | 10.0 | 20.0 | 50.0 | 100 |
| Ethyl acetate | FB | Ave | 61731<br>2713629 | 116657 | 235203 | 666637 | 1307022 | 10.0<br>400 | 20.0 | 40.0 | 100 | 200 |
| Propionitrile | FB | Ave | 52860<br>2001714 | 94565 | 174727 | 500380 | 942722 | 50.0<br>2000 | 100 | 200 | 500 | 1000 |
| Methacrylonitrile | FB | Ave | 264644<br>10194839 | 475891 | 913698 | 2507397 | 4932151 | 50.0<br>2000 | 100 | 200 | 500 | 1000 |
| Tert-amyl methyl ether | FB | Ave | 37226<br>1851023 | 71362 | 144295 | 414571 | 874489 | 5.00<br>200 | 10.0 | 20.0 | 50.0 | 100 |
| n-Butanol | FB | Qua | 18792<br>784490 | 28157 | 45708 | 172807 | 331694 | 125<br>5000 | 250 | 500 | 1250 | 2500 |
| Ethyl acrylate | FB | Ave | 30789<br>1547078 | 58546 | 130182 | 369840 | 770109 | 5.00<br>200 | 10.0 | 20.0 | 50.0 | 100 |
| Methyl methacrylate | FB | Ave | 60925<br>2691827 | 116812 | 237896 | 655501 | 1360547 | 10.0<br>400 | 20.0 | 40.0 | 100 | 200 |
| 2-Nitropropane | FB | Qua | 9958<br>744616 | 20703 | 40995 | 150340 | 323825 | 10.0<br>400 | 20.0 | 40.0 | 100 | 200 |
| n-Butyl acetate | FB | Ave | 51170<br>1926459 | 89406 | 164263 | 487893 | 1001885 | 5.00<br>200 | 10.0 | 20.0 | 50.0 | 100 |
| 1-Chlorohexane | CBZ | Ave | 25554<br>1065503 | 48122 | 99357 | 244163 | 549387 | 5.00<br>200 | 10.0 | 20.0 | 50.0 | 100 |
| Cyclohexanone | CBZ | Qua | 14662<br>433012 | 22018 | 33024 | 108174 | 196967 | 50.0<br>2000 | 100 | 200 | 500 | 1000 |
| 1,2,3-Trimethylbenzene | DCB | Ave | 99284<br>4034098 | 189500 | 360620 | 942781 | 2000672 | 5.00<br>200 | 10.0 | 20.0 | 50.0 | 100 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 189147

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/15/2015 01:51 Calibration End Date: 07/15/2015 03:38 Calibration ID: 29499

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| 1,3,5-Trichlorobenzene | DCB | Ave | 41531<br>1610278 | 78143 | 150029 | 380909 | 788451 | 5.00<br>200 | 10.0 | 20.0 | 50.0 | 100 |
| 2-Methylnaphthalene | DCB | Ave | 135125<br>3624085 | 228819 | 431799 | 1104036 | 2151886 | 10.0<br>400 | 20.0 | 40.0 | 100 | 200 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Qua = Quadratic ISTD |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35 Calibration End Date: 07/30/2015 05:04 Calibration ID: 29903

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-191232/10 | UX88725.D |
| Level 2 | STD8260 240-191232/9 | UX88724.D |
| Level 3 | STD8260 240-191232/8 | UX88723.D |
| Level 4 | STD8260 240-191232/7 | UX88722.D |
| Level 5 | ICIS 240-191232/6 | UX88721.D |
| Level 6 | STD8260 240-191232/5 | UX88720.D |
| Level 7 | STD8260 240-191232/4 | UX88719.D |
| Level 8 | STD8260 240-191232/3 | UX88718.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | 0.1321<br>0.1303 | 0.1134<br>0.1532 | 0.1571<br>0.1672 | 0.1280 | 0.1474 | Ave | | 0.1411 | | | 0.1000 | 12.7 | | 15.0 | | | |
| Chloromethane | 0.2636<br>0.2350 | 0.2111<br>0.2667 | 0.2993<br>0.2673 | 0.2283 | 0.2646 | Ave | | 0.2545 | | | 0.1000 | 11.0 | | 15.0 | | | |
| Vinyl chloride | 0.1348<br>0.1634 | 0.1450<br>0.1856 | 0.2126<br>0.1909 | 0.1655 | 0.1888 | Ave | | 0.1733 | | | 0.1000 | 14.9 | | 15.0 | | | |
| Butadiene | 0.2213<br>0.2122 | 0.2009<br>0.2434 | 0.2963<br>0.2578 | 0.2093 | 0.2447 | Ave | | 0.2357 | | | | 13.4 | | 15.0 | | | |
| Bromomethane | 0.1060<br>0.0766 | 0.0664<br>0.0743 | 0.1106<br>0.0718 | 0.0793 | 0.0925 | Qua | 0.0948 | 0.0765 | -0.000026 | | 0.0500 | | | | 1.0000 | | 0.9900 |
| Chloroethane | +++++<br>0.1175 | +++++<br>0.1203 | 0.1599<br>0.1122 | 0.1095 | 0.1333 | Ave | | 0.1254 | | | | 15.0 | | 15.0 | | | |
| Dichlorofluoromethane | 0.2789<br>0.2670 | 0.2303<br>0.2863 | 0.3705<br>0.2872 | 0.2676 | 0.2967 | Ave | | 0.2856 | | | | 14.0 | | 15.0 | | | |
| Trichlorofluoromethane | +++++<br>0.2816 | 0.2474<br>0.3037 | 0.3313<br>0.3036 | 0.2858 | 0.2937 | Ave | | 0.2924 | | | 0.1000 | 8.8 | | 15.0 | | | |
| Ethyl ether | 0.3290<br>0.2911 | 0.2844<br>0.2745 | 0.3352<br>0.3064 | 0.2732 | 0.2945 | Ave | | 0.2985 | | | | 7.8 | | 15.0 | | | |
| Acrolein | 0.0401<br>0.0299 | 0.0285<br>0.0319 | 0.0340<br>0.0322 | 0.0291 | 0.0339 | Ave | | 0.0324 | | | | 11.5 | | 15.0 | | | |
| 1,1-Dichloroethene | +++++<br>0.2263 | 0.1859<br>0.2212 | 0.2757<br>0.2397 | 0.1969 | 0.2287 | Ave | | 0.2249 | | | 0.1000 | 13.0 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | +++++<br>0.1137 | 0.0845<br>0.1189 | 0.1311<br>0.1255 | 0.0964 | 0.1209 | Ave | | 0.1130 | | | 0.0500 | 14.8 | | 15.0 | | | |
| Acetone | +++++<br>0.1711 | 0.3313<br>0.1720 | 0.2706<br>0.1655 | 0.1948 | 0.2037 | Lin1 | 0.7884 | 0.1657 | | | 0.0500 | | | | 0.9980 | | 0.9900 |
| Iodomethane | 0.3851<br>0.3876 | 0.3238<br>0.3697 | 0.4722<br>0.3893 | 0.3363 | 0.3801 | Ave | | 0.3805 | | | | 11.7 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35  Calibration End Date: 07/30/2015 05:04  Calibration ID: 29903

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon disulfide | +++++<br>0.8290 | 0.6093<br>0.8168 | 0.8467<br>0.8950 | 0.6645 | 0.7915 | Ave | | 0.7790 | | | 0.1000 | 13.3 | | 15.0 | | | |
| 3-Chloro-1-propene | 0.1207<br>0.1502 | 0.1138<br>0.1482 | 0.1516<br>0.1562 | 0.1270 | 0.1511 | Ave | | 0.1399 | | | | 11.8 | | 15.0 | | | |
| Methyl acetate | 0.3354<br>0.3142 | 0.3163<br>0.3301 | 0.3511<br>0.3448 | 0.3155 | 0.3391 | Ave | | 0.3308 | | | 0.1000 | 4.3 | | 15.0 | | | |
| Methylene Chloride | +++++<br>0.3064 | +++++<br>0.2798 | 0.5985<br>0.2919 | 0.3581 | 0.3500 | Lin1 | 1.3334 | 0.2785 | | | 0.1000 | | | | 0.9970 | | 0.9900 |
| 2-Methyl-2-propanol | +++++<br>0.0277 | +++++<br>0.0272 | 0.0338<br>0.0245 | 0.0235 | 0.0306 | Ave | | 0.0279 | | | | 13.8 | | 15.0 | | | |
| Acrylonitrile | 0.1583<br>0.1377 | 0.1330<br>0.1420 | 0.1548<br>0.1447 | 0.1315 | 0.1480 | Ave | | 0.1437 | | | | 6.7 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.8187<br>0.8349 | 0.7042<br>0.8119 | 0.9030<br>0.8342 | 0.7217 | 0.8500 | Ave | | 0.8098 | | | 0.1000 | 8.2 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.2329<br>0.2670 | 0.2115<br>0.2517 | 0.3039<br>0.2653 | 0.2253 | 0.2651 | Ave | | 0.2529 | | | 0.1000 | 11.6 | | 15.0 | | | |
| Hexane | 0.0890<br>0.0735 | 0.0862<br>0.0721 | 0.0932<br>0.0746 | 0.0702 | 0.0686 | Ave | | 0.0784 | | | | 12.1 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.5574<br>0.5921 | 0.5187<br>0.5725 | 0.6500<br>0.6103 | 0.5359 | 0.5798 | Ave | | 0.5771 | | | 0.1000 | 7.2 | | 15.0 | | | |
| Vinyl acetate | 0.4178<br>0.5446 | 0.3761<br>0.5673 | 0.4691<br>0.4212 | 0.4118 | 0.4997 | Ave | | 0.4635 | | | | 14.8 | | 15.0 | | | |
| 2-Butanone (MEK) | 0.2390<br>0.2004 | 0.2440<br>0.2082 | 0.2257<br>0.1964 | 0.2198 | 0.2153 | Ave | | 0.2186 | | | 0.0500 | 7.8 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.2755<br>0.2800 | 0.2525<br>0.2665 | 0.3252<br>0.2808 | 0.2558 | 0.2769 | Ave | | 0.2767 | | | 0.1000 | 8.1 | | 15.0 | | | |
| 2,2-Dichloropropane | 0.1634<br>0.2832 | 0.1626<br>0.2819 | 0.2377<br>0.3000 | 0.1911 | 0.2490 | Lin1 | -0.243 | 0.2909 | | | | | | | 0.9960 | | 0.9900 |
| Chlorobromomethane | 0.1008<br>0.1222 | 0.1092<br>0.1181 | 0.1451<br>0.1235 | 0.1139 | 0.1240 | Ave | | 0.1196 | | | | 10.9 | | 15.0 | | | |
| Tetrahydrofuran | 0.1347<br>0.1229 | 0.1241<br>0.1272 | 0.1289<br>0.1259 | 0.1244 | 0.1330 | Ave | | 0.1276 | | | | 3.4 | | 15.0 | | | |
| Chloroform | 0.4864<br>0.4860 | 0.4514<br>0.4611 | 0.5409<br>0.4824 | 0.4469 | 0.4740 | Ave | | 0.4786 | | | | 6.2 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.2548<br>0.3771 | 0.2875<br>0.3627 | 0.3746<br>0.3858 | 0.3120 | 0.3548 | Ave | | 0.3387 | | | 0.1000 | 14.2 | | 15.0 | | | |
| Cyclohexane | +++++<br>0.6285 | 0.5035<br>0.6114 | 0.6700<br>0.6338 | 0.5607 | 0.6245 | Ave | | 0.6046 | | | 0.1000 | 9.1 | | 15.0 | | | |
| 1,1-Dichloropropene | 0.2436<br>0.3492 | 0.2972<br>0.3396 | 0.3658<br>0.3545 | 0.3310 | 0.3349 | Ave | | 0.3270 | | | | 12.0 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35  Calibration End Date: 07/30/2015 05:04  Calibration ID: 29903

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon tetrachloride | +++++<br>0.3150 | +++++<br>0.3123 | 0.2806<br>0.3387 | 0.2486 | 0.2826 | Ave | | 0.2963 | | | 0.1000 | 10.8 | | 15.0 | | | |
| Isobutyl alcohol | 0.0150<br>0.0120 | 0.0102<br>0.0119 | 0.0149<br>0.0105 | 0.0117 | 0.0129 | Ave | | 0.0124 | | | | 14.4 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.4878<br>0.4489 | 0.4876<br>0.4331 | 0.4971<br>0.4344 | 0.4619 | 0.4470 | Ave | | 0.4622 | | | 0.1000 | 5.5 | | 15.0 | | | |
| Benzene | 1.0013<br>1.0551 | 1.0073<br>1.0086 | 1.1409<br>1.0494 | 0.9988 | 1.0195 | Ave | | 1.0351 | | | 0.5000 | 4.6 | | 15.0 | | | |
| n-Heptane | +++++<br>0.0718 | 0.0621<br>0.0705 | 0.0775<br>0.0718 | 0.0690 | 0.0668 | Ave | | 0.0699 | | | | 6.8 | | 15.0 | | | |
| Trichloroethene | 0.2239<br>0.2327 | 0.2323<br>0.2307 | 0.2581<br>0.2415 | 0.2354 | 0.2254 | Ave | | 0.2350 | | | 0.2000 | 4.6 | | 15.0 | | | |
| Methylcyclohexane | +++++<br>0.4707 | 0.3791<br>0.4551 | 0.5020<br>0.4852 | 0.4076 | 0.4560 | Ave | | 0.4508 | | | 0.1000 | 9.6 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.2306<br>0.2906 | 0.2830<br>0.2784 | 0.2849<br>0.2827 | 0.2866 | 0.2757 | Ave | | 0.2766 | | | 0.1000 | 6.9 | | 15.0 | | | |
| 1,4-Dioxane | 0.0022<br>0.0022 | 0.0019<br>0.0020 | 0.0026<br>0.0019 | 0.0021 | 0.0023 | Ave | | 0.0021 | | | | 10.7 | | 15.0 | | | |
| Dibromomethane | 0.1569<br>0.1384 | 0.1424<br>0.1365 | 0.1563<br>0.1361 | 0.1345 | 0.1351 | Ave | | 0.1420 | | | | 6.6 | | 15.0 | | | |
| Dichlorobromomethane | 0.2676<br>0.3273 | 0.2944<br>0.3223 | 0.2959<br>0.3330 | 0.2987 | 0.3002 | Ave | | 0.3049 | | | 0.1000 | 7.0 | | 15.0 | | | |
| 2-Chloroethyl vinyl ether | +++++<br>0.0080 | +++++<br>0.0100 | 0.0046<br>0.0099 | 0.0085 | 0.0062 | Lin1 | -0.062 | 0.0098 | | | | | | | 0.9920 | | 0.9900 |
| cis-1,3-Dichloropropene | +++++<br>0.3370 | 0.2554<br>0.3442 | 0.2510<br>0.3511 | 0.3071 | 0.2799 | Ave | | 0.3036 | | | 0.1500 | 13.9 | | 15.0 | | | |
| 4-Methyl-2-pentanone (MIBK) | 0.3441<br>0.3695 | 0.3474<br>0.3838 | 0.3633<br>0.3520 | 0.3808 | 0.3788 | Ave | | 0.3650 | | | 0.0500 | 4.3 | | 15.0 | | | |
| Toluene | 1.5154<br>1.5909 | 1.5612<br>1.5632 | 1.6996<br>1.6699 | 1.5226 | 1.5117 | Ave | | 1.5793 | | | 0.4000 | 4.5 | | 15.0 | | | |
| trans-1,3-Dichloropropene | 0.2045<br>0.4159 | 0.2604<br>0.4511 | 0.2731<br>0.4740 | 0.3419 | 0.3375 | Qua | -0.783 | 0.4260 | 0.0002631 | | 0.1000 | | | | 1.0000 | | 0.9900 |
| Ethyl methacrylate | +++++<br>0.4407 | 0.3513<br>0.4554 | 0.3902<br>0.4406 | 0.4361 | 0.4254 | Ave | | 0.4200 | | | | 8.7 | | 15.0 | | | |
| 1,1,2-Trichloroethane | 0.2435<br>0.2780 | 0.3132<br>0.2804 | 0.3064<br>0.2816 | 0.2922 | 0.2709 | Ave | | 0.2833 | | | 0.1000 | 7.6 | | 15.0 | | | |
| 1,3-Dichloropropane | 0.4856<br>0.5423 | 0.5703<br>0.5528 | 0.5925<br>0.5461 | 0.5803 | 0.5378 | Ave | | 0.5510 | | | | 5.9 | | 15.0 | | | |
| Tetrachloroethene | 0.2330<br>0.3023 | 0.2810<br>0.3014 | 0.3263<br>0.3151 | 0.2940 | 0.3023 | Ave | | 0.2944 | | | 0.2000 | 9.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35  Calibration End Date: 07/30/2015 05:04  Calibration ID: 29903

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2-Hexanone | 0.4589<br>0.3793 | 0.3890<br>0.4006 | 0.4031<br>0.3485 | 0.4264 | 0.3942 | Ave | | 0.4000 | | | 0.0500 | 8.1 | | 15.0 | | | |
| Chlorodibromomethane | +++++<br>0.2915 | 0.2138<br>0.2943 | 0.2332<br>0.3111 | 0.2507 | 0.2583 | Ave | | 0.2647 | | | | 13.4 | | 15.0 | | | |
| Ethylene Dibromide | 0.2347<br>0.2762 | 0.2906<br>0.2856 | 0.2933<br>0.2798 | 0.2749 | 0.2702 | Ave | | 0.2757 | | | | 6.7 | | 15.0 | | | |
| Chlorobenzene | 0.8857<br>0.9308 | 0.9993<br>0.9072 | 1.0209<br>0.9233 | 0.9484 | 0.8972 | Ave | | 0.9391 | | | 0.3000 | 5.1 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.2603<br>0.3362 | 0.2227<br>0.3249 | 0.3079<br>0.3316 | 0.2611 | 0.3075 | Ave | | 0.2940 | | | | 14.0 | | 15.0 | | | |
| Ethylbenzene | 0.5325<br>0.5156 | 0.5153<br>0.4931 | 0.5699<br>0.4935 | 0.5118 | 0.5023 | Ave | | 0.5167 | | | | 4.9 | | 15.0 | | | |
| m-Xylene & p-Xylene | 0.5111<br>0.6338 | 0.5640<br>0.5968 | 0.7337<br>0.5975 | 0.6178 | 0.6263 | Ave | | 0.6101 | | | | 10.4 | | 15.0 | | | |
| o-Xylene | 0.5860<br>0.6234 | 0.5597<br>0.5818 | 0.7228<br>0.5693 | 0.5925 | 0.6323 | Ave | | 0.6085 | | | | 8.6 | | 15.0 | | | |
| Styrene | 0.8663<br>0.9946 | 0.9107<br>0.9432 | 1.0105<br>0.9228 | 0.9977 | 0.9613 | Ave | | 0.9509 | | | 0.3000 | 5.2 | | 15.0 | | | |
| Bromoform | +++++<br>0.1688 | +++++<br>0.1725 | 0.1346<br>0.1745 | 0.1341 | 0.1484 | Ave | | 0.1555 | | | 0.1000 | 12.1 | | 15.0 | | | |
| Isopropylbenzene | 1.2399<br>1.6788 | 1.3621<br>1.5275 | 1.7949<br>1.5054 | 1.5320 | 1.6682 | Ave | | 1.5386 | | | 0.1000 | 11.6 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 1.1178<br>0.8460 | 0.7769<br>0.8566 | 0.9350<br>0.7781 | 0.8033 | 0.8300 | Ave | | 0.8680 | | | 0.3000 | 13.0 | | 15.0 | | | |
| Bromobenzene | 0.8849<br>0.9203 | 1.0157<br>0.8817 | 0.9194<br>0.8479 | 0.9962 | 0.8191 | Ave | | 0.9106 | | | | 7.5 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.3645<br>0.2759 | 0.2752<br>0.2852 | 0.3450<br>0.2570 | 0.3015 | 0.2738 | Ave | | 0.2973 | | | | 12.8 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | +++++<br>0.2732 | 0.2343<br>0.3026 | 0.2506<br>0.2752 | 0.2666 | 0.2431 | Ave | | 0.2637 | | | | 8.8 | | 15.0 | | | |
| N-Propylbenzene | 0.6914<br>1.0471 | 0.9318<br>0.9818 | 0.9827<br>0.9515 | 1.0468 | 0.9593 | Ave | | 0.9490 | | | | 11.8 | | 15.0 | | | |
| 2-Chlorotoluene | 0.6794<br>0.8572 | 0.7938<br>0.8212 | 0.9496<br>0.7925 | 0.8401 | 0.7912 | Ave | | 0.8156 | | | | 9.3 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 2.1891<br>3.3370 | 2.9109<br>3.1719 | 3.3711<br>3.1546 | 3.1409 | 3.0793 | Ave | | 3.0443 | | | | 12.3 | | 15.0 | | | |
| 4-Chlorotoluene | 0.8045<br>0.8207 | 0.8195<br>0.7984 | 0.8281<br>0.7676 | 0.8887 | 0.7809 | Ave | | 0.8136 | | | | 4.5 | | 15.0 | | | |
| tert-Butylbenzene | 1.7942<br>2.8350 | 2.3516<br>2.6952 | 2.5292<br>2.9278 | 2.7401 | 2.5233 | Ave | | 2.5495 | | | | 14.0 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35 Calibration End Date: 07/30/2015 05:04 Calibration ID: 29903

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,2,4-Trimethylbenzene | 2.4658<br>3.2632 | 2.8912<br>3.1305 | 3.4321<br>3.1783 | 3.1148 | 3.1141 | Ave | | 3.0738 | | | | 9.4 | | 15.0 | | | |
| sec-Butylbenzene | 2.4505<br>3.9697 | 3.3958<br>3.8627 | 3.9265<br>3.9512 | 3.7820 | 3.7587 | Ave | | 3.6371 | | | | 14.1 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.6380<br>1.4856 | 1.6175<br>1.5113 | 1.7985<br>1.4747 | 1.5773 | 1.4995 | Ave | | 1.5753 | | | 0.6000 | 6.9 | | 15.0 | | | |
| 4-Isopropyltoluene | 2.0429<br>3.0339 | 2.6459<br>2.9986 | 3.1244<br>3.1413 | 2.8616 | 2.9162 | Ave | | 2.8456 | | | | 12.7 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 1.7161<br>1.4483 | 1.6014<br>1.4899 | 1.7260<br>1.4814 | 1.5232 | 1.4329 | Ave | | 1.5524 | | | 0.5000 | 7.5 | | 15.0 | | | |
| n-Butylbenzene | +++++<br>2.6320 | 2.0158<br>2.6928 | 2.8623<br>2.8570 | 2.4030 | 2.5949 | Ave | | 2.5797 | | | | 11.4 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.9126<br>1.3749 | 1.4808<br>1.4460 | 1.7342<br>1.4671 | 1.4089 | 1.4240 | Ave | | 1.5311 | | | 0.4000 | 12.4 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | +++++<br>0.1259 | +++++<br>0.1571 | 0.1140<br>0.1755 | 0.0955 | 0.1183 | Qua | -0.401 | 0.1355 | 0.0002121 | | 0.0500 | | | | 0.9990 | | 0.9900 |
| 1,2,4-Trichlorobenzene | 1.2873<br>1.0593 | 0.9453<br>1.1962 | 1.4224<br>1.3439 | 0.8759 | 1.1578 | Lin1 | -0.397 | 1.2545 | | | 0.2000 | | | | 0.9900 | | 0.9900 |
| Hexachlorobutadiene | +++++<br>0.6503 | +++++<br>0.7088 | 0.7353<br>0.7901 | 0.5587 | 0.6655 | Ave | | 0.6848 | | | | 11.6 | | 15.0 | | | |
| Naphthalene | 3.5776<br>2.2692 | 1.9057<br>2.6058 | 2.8701<br>2.8241 | 1.8628 | 2.5181 | Qua | -0.246 | 2.2907 | 0.0026922 | | | | | | 1.0000 | | 0.9900 |
| 1,2,3-Trichlorobenzene | 1.7206<br>1.0468 | 0.9822<br>1.1480 | 1.4853<br>1.2586 | 0.8691 | 1.1572 | Lin1 | 0.0849 | 1.1884 | | | | | | | 0.9920 | | 0.9900 |
| Dibromofluoromethane (Surr) | +++++<br>0.2299 | +++++<br>0.2198 | 0.5882<br>0.2311 | 0.3575 | 0.2798 | Lin1 | 1.6313 | 0.2130 | | | | | | | 0.9950 | | 0.9900 |
| 1,2-Dichloroethane-d4 (Surr) | +++++<br>0.3117 | +++++<br>0.3245 | 0.8497<br>0.3370 | 0.5850 | 0.4251 | Lin1 | 2.5528 | 0.3083 | | | | | | | 0.9930 | | 0.9900 |
| Toluene-d8 (Surr) | +++++<br>1.3034 | +++++<br>1.2868 | 3.2443<br>1.3682 | 2.1008 | 1.5924 | Lin1 | 8.6927 | 1.2582 | | | | | | | 0.9940 | | 0.9900 |
| 4-Bromofluorobenzene (Surr) | +++++<br>0.3937 | +++++<br>0.3830 | 1.0904<br>0.3654 | 0.7675 | 0.5368 | Lin1 | 3.8505 | 0.3424 | | | | | | | 0.9980 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35 Calibration End Date: 07/30/2015 05:04 Calibration ID: 29903

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-191232/10 | UX88725.D |
| Level 2 | STD8260 240-191232/9 | UX88724.D |
| Level 3 | STD8260 240-191232/8 | UX88723.D |
| Level 4 | STD8260 240-191232/7 | UX88722.D |
| Level 5 | ICIS 240-191232/6 | UX88721.D |
| Level 6 | STD8260 240-191232/5 | UX88720.D |
| Level 7 | STD8260 240-191232/4 | UX88719.D |
| Level 8 | STD8260 240-191232/3 | UX88718.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1 LVL 6 | LVL 2 LVL 7 | LVL 3 LVL 8 | LVL 4 | LVL 5 | LVL 1 LVL 6 | LVL 2 LVL 7 | LVL 3 LVL 8 | LVL 4 | LVL 5 |
| Dichlorodifluoromethane | FB | Ave | 3474<br>215398 | 8087<br>475117 | 20438<br>1007796 | 46757 | 82812 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Chloromethane | FB | Ave | 6930<br>388593 | 15051<br>826915 | 38955<br>1611037 | 83414 | 148640 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Vinyl chloride | FB | Ave | 3544<br>270172 | 10339<br>575508 | 27670<br>1150663 | 60466 | 106043 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Butadiene | FB | Ave | 5819<br>350930 | 14326<br>754819 | 38557<br>1553429 | 76475 | 137459 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Bromomethane | FB | Qua | 2787<br>126612 | 4734<br>230400 | 14393<br>432898 | 28988 | 51977 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Chloroethane | FB | Ave | +++++<br>194201 | +++++<br>372932 | 20803<br>676029 | 40026 | 74866 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Dichlorofluoromethane | FB | Ave | 7332<br>441555 | 16417<br>887850 | 48218<br>1730860 | 97788 | 166694 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Trichlorofluoromethane | FB | Ave | +++++<br>465612 | 17638<br>941746 | 43115<br>1829523 | 104425 | 165026 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Ethyl ether | FB | Ave | 8650<br>481320 | 20275<br>851240 | 43627<br>1846389 | 99831 | 165442 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Acrolein | FB | Ave | 5276<br>246804 | 10142<br>493854 | 22099<br>969388 | 53238 | 95240 | 5.00<br>250 | 10.0<br>500 | 25.0<br>1000 | 50.0 | 100 |
| 1,1-Dichloroethene | FB | Ave | +++++<br>374206 | 13255<br>686008 | 35873<br>1444827 | 71946 | 128504 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | FB | Ave | +++++<br>188054 | 6022<br>368648 | 17065<br>756434 | 35213 | 67913 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Acetone | FB | Lin1 | +++++<br>565690 | 47240<br>1066719 | 70438<br>1995245 | 142344 | 228883 | +++++<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| Iodomethane | FB | Ave | 10124<br>640930 | 23087<br>1146413 | 61451<br>2346077 | 122890 | 213572 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Carbon disulfide | FB | Ave | +++++<br>1370798 | 43441<br>2532638 | 110188<br>5393852 | 242831 | 444651 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35 Calibration End Date: 07/30/2015 05:04 Calibration ID: 29903

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 |
| 3-Chloro-1-propene | FB | Ave | 3174<br>248371 | 8111<br>459550 | 19732<br>941156 | 46426 | 84888 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Methyl acetate | FB | Ave | 44082<br>2597514 | 112759<br>5117451 | 228422<br>10390579 | 576496 | 952461 | 5.00<br>250 | 10.0<br>500 | 25.0<br>1000 | 50.0 | 100 |
| Methylene Chloride | FB | Lin1 | +++++<br>506547 | +++++<br>867437 | 77886<br>1758980 | 130876 | 196633 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2-Methyl-2-propanol | FB | Ave | +++++<br>458456 | +++++<br>841872 | 44023<br>1477308 | 85925 | 171659 | +++++<br>500 | +++++<br>1000 | 50.0<br>2000 | 100 | 200 |
| Acrylonitrile | FB | Ave | 41604<br>2276996 | 94854<br>4402094 | 201432<br>8719139 | 480494 | 831302 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100 | 200 |
| Methyl tert-butyl ether | FB | Ave | 21524<br>1380396 | 50202<br>2517525 | 117506<br>5027382 | 263716 | 477515 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| trans-1,2-Dichloroethene | FB | Ave | 6123<br>441489 | 15081<br>780372 | 39554<br>1599085 | 82325 | 148960 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Hexane | FB | Ave | 2340<br>121473 | 6149<br>223660 | 12125<br>449870 | 25657 | 38539 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1-Dichloroethane | FB | Ave | 14653<br>978972 | 36980<br>1775266 | 84582<br>3678291 | 195825 | 325730 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Vinyl acetate | FB | Ave | 10984<br>900552 | 26810<br>1758987 | 61044<br>2538569 | 150494 | 280750 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2-Butanone (MEK) | FB | Ave | 12566<br>662691 | 34797<br>1291450 | 58740<br>2367067 | 160626 | 241864 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| cis-1,2-Dichloroethene | FB | Ave | 7244<br>462961 | 17999<br>826232 | 42321<br>1692517 | 93478 | 155557 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2,2-Dichloropropane | FB | Lin1 | 4296<br>468274 | 11593<br>874180 | 30936<br>1808143 | 69826 | 139865 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Chlorobromomethane | FB | Ave | 2649<br>202074 | 7783<br>366251 | 18888<br>744297 | 41627 | 69660 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Tetrahydrofuran | FB | Ave | 7084<br>406513 | 17688<br>788681 | 33544<br>1518106 | 90942 | 149470 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| Chloroform | FB | Ave | 12788<br>803613 | 32180<br>1429887 | 70395<br>2907092 | 163299 | 266320 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1,1-Trichloroethane | FB | Ave | 6699<br>623476 | 20498<br>1124791 | 48752<br>2325338 | 114020 | 199341 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Cyclohexane | FB | Ave | +++++<br>1039262 | 35895<br>1895812 | 87191<br>3819980 | 204896 | 350853 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1-Dichloropropene | FB | Ave | 6403<br>577421 | 21185<br>1053095 | 47605<br>2136628 | 120951 | 188172 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Carbon tetrachloride | FB | Ave | +++++<br>520790 | +++++<br>968265 | 36521<br>2041474 | 90834 | 158767 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Isobutyl alcohol | FB | Ave | 9845<br>496725 | 18100<br>922435 | 48378<br>1586721 | 107293 | 181553 | 25.0<br>1250 | 50.0<br>2500 | 125<br>5000 | 250 | 500 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35 Calibration End Date: 07/30/2015 05:04 Calibration ID: 29903

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 |
| 1,2-Dichloroethane | FB | Ave | 12825<br>742241 | 34763<br>1343026 | 64694<br>2618263 | 168804 | 251137 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Benzene | FB | Ave | 26325<br>1744546 | 71817<br>3127556 | 148470<br>6324576 | 365003 | 572779 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| n-Heptane | FB | Ave | +++++<br>118695 | 4427<br>218752 | 10090<br>432763 | 25226 | 37554 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Trichloroethene | FB | Ave | 5887<br>384809 | 16560<br>715274 | 33589<br>1455286 | 86037 | 126619 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Methylcyclohexane | FB | Ave | +++++<br>778240 | 27025<br>1411203 | 65331<br>2923845 | 148967 | 256199 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2-Dichloropropane | FB | Ave | 6063<br>480527 | 20173<br>863200 | 37078<br>1703845 | 104723 | 154868 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,4-Dioxane | FB | Ave | 1147<br>71927 | 2689<br>121705 | 6670<br>227257 | 15559 | 25856 | 20.0<br>1000 | 40.0<br>2000 | 100<br>4000 | 200 | 400 |
| Dibromomethane | FB | Ave | 4125<br>228773 | 10155<br>423286 | 20342<br>820019 | 49148 | 75882 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Dichlorobromomethane | FB | Ave | 7036<br>541112 | 20989<br>999368 | 38506<br>2006716 | 109160 | 168635 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2-Chloroethyl vinyl ether | FB | Lin1 | +++++<br>26619 | +++++<br>61713 | 1191<br>118924 | 6248 | 6945 | +++++<br>100 | +++++<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| cis-1,3-Dichloropropene | FB | Ave | +++++<br>557159 | 18205<br>1067131 | 32664<br>2115778 | 112216 | 157235 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | FB | Ave | 18095<br>1221930 | 49532<br>2380397 | 94566<br>4242931 | 278306 | 425603 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| Toluene | CBZ | Ave | 22064<br>1551933 | 70621<br>2781164 | 122902<br>5466002 | 353822 | 482509 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| trans-1,3-Dichloropropene | CBZ | Qua | 2977<br>405684 | 11781<br>802553 | 19747<br>1551499 | 79460 | 107712 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Ethyl methacrylate | CBZ | Ave | +++++<br>429950 | 15891<br>810182 | 28220<br>1442115 | 101344 | 135771 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1,2-Trichloroethane | CBZ | Ave | 3546<br>271178 | 14166<br>498965 | 22155<br>921874 | 67905 | 86457 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,3-Dichloropropane | CBZ | Ave | 7071<br>529038 | 25799<br>983577 | 42844<br>1787500 | 134844 | 171656 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Tetrachloroethene | CBZ | Ave | 3392<br>294930 | 12713<br>536185 | 23599<br>1031265 | 68313 | 96485 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2-Hexanone | CBZ | Ave | 13362<br>739956 | 35190<br>1425410 | 58299<br>2281643 | 198186 | 251682 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| Chlorodibromomethane | CBZ | Ave | +++++<br>284330 | 9669<br>523664 | 16863<br>1018379 | 58266 | 82461 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Ethylene Dibromide | CBZ | Ave | 3417<br>269468 | 13147<br>508160 | 21211<br>915674 | 63877 | 86243 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35 Calibration End Date: 07/30/2015 05:04 Calibration ID: 29903

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 |
| Chlorobenzene | CBZ | Ave | 12896<br>908035 | 45204<br>1614116 | 73822<br>3021966 | 220385 | 286379 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1,1,2-Tetrachloroethane | CBZ | Ave | 3790<br>327964 | 10075<br>578063 | 22263<br>1085341 | 60676 | 98153 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Ethylbenzene | CBZ | Ave | 7753<br>502949 | 23309<br>877372 | 41208<br>1615386 | 118923 | 160342 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| m-Xylene & p-Xylene | CBZ | Ave | 7442<br>618258 | 25510<br>1061760 | 53053<br>1955816 | 143563 | 199922 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| o-Xylene | CBZ | Ave | 8532<br>608087 | 25316<br>1035179 | 52271<br>1863497 | 137690 | 201835 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Styrene | CBZ | Ave | 12613<br>970267 | 41194<br>1678160 | 73072<br>3020605 | 231838 | 306845 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Bromoform | CBZ | Ave | +++++<br>164641 | +++++<br>306960 | 9734<br>571250 | 31154 | 47358 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Isopropylbenzene | CBZ | Ave | 18054<br>1637670 | 61614<br>2717802 | 129792<br>4927519 | 355999 | 532483 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1,2,2-Tetrachloroethane | DCB | Ave | 6700<br>296525 | 11939<br>533483 | 27248<br>891333 | 65067 | 105628 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Bromobenzene | DCB | Ave | 5304<br>322557 | 15609<br>549088 | 26793<br>971254 | 80691 | 104239 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2,3-Trichloropropane | DCB | Ave | 2185<br>96692 | 4229<br>177618 | 10055<br>294392 | 24421 | 34840 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| trans-1,4-Dichloro-2-butene | DCB | Ave | +++++<br>95739 | 3600<br>188429 | 7302<br>315306 | 21597 | 30943 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| N-Propylbenzene | DCB | Ave | 4144<br>366998 | 14319<br>611453 | 28637<br>1089957 | 84787 | 122083 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2-Chlorotoluene | DCB | Ave | 4072<br>300431 | 12199<br>511446 | 27674<br>907841 | 68050 | 100686 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,3,5-Trimethylbenzene | DCB | Ave | 13121<br>1169588 | 44732<br>1975399 | 98237<br>3613779 | 254411 | 391875 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 4-Chlorotoluene | DCB | Ave | 4822<br>287644 | 12593<br>497220 | 24131<br>879335 | 71988 | 99384 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| tert-Butylbenzene | DCB | Ave | 10754<br>993643 | 36137<br>1678494 | 73704<br>3353932 | 221951 | 321117 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2,4-Trimethylbenzene | DCB | Ave | 14780<br>1143743 | 44429<br>1949583 | 100016<br>3640920 | 252297 | 396302 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| sec-Butylbenzene | DCB | Ave | 14688<br>1391355 | 52184<br>2405595 | 114424<br>4526241 | 306344 | 478337 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,3-Dichlorobenzene | DCB | Ave | 9818<br>520708 | 24857<br>941170 | 52412<br>1689365 | 127765 | 190835 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 4-Isopropyltoluene | DCB | Ave | 12245<br>1063373 | 40660<br>1867477 | 91050<br>3598534 | 231794 | 371116 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 191232

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 07/30/2015 02:35 Calibration End Date: 07/30/2015 05:04 Calibration ID: 29903

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 |
| 1,4-Dichlorobenzene | DCB | Ave | 10286<br>507629 | 24609<br>927871 | 50299<br>1697055 | 123381 | 182351 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| n-Butylbenzene | DCB | Ave | +++++<br>922495 | 30977<br>1677021 | 83412<br>3272827 | 194646 | 330234 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2-Dichlorobenzene | DCB | Ave | 11464<br>481899 | 22755<br>900531 | 50538<br>1680638 | 114123 | 181216 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | DCB | Qua | +++++<br>44124 | +++++<br>97839 | 3321<br>201076 | 7732 | 15050 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2,4-Trichlorobenzene | DCB | Lin1 | 7716<br>371288 | 14526<br>744987 | 41451<br>1539552 | 70950 | 147347 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Hexachlorobutadiene | DCB | Ave | +++++<br>227915 | +++++<br>441429 | 21427<br>905098 | 45253 | 84698 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Naphthalene | DCB | Qua | 21444<br>795334 | 29285<br>1622846 | 83637<br>3235092 | 150891 | 320459 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2,3-Trichlorobenzene | DCB | Lin1 | 10313<br>366883 | 15094<br>714939 | 43284<br>1441752 | 70395 | 147264 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Dibromofluoromethane (Surr) | FB | Lin1 | +++++<br>380111 | +++++<br>681545 | 76539<br>1392712 | 130630 | 157205 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Lin1 | +++++<br>515428 | +++++<br>1006250 | 110575<br>2030768 | 213759 | 238799 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Toluene-d8 (Surr) | CBZ | Lin1 | +++++<br>1271417 | +++++<br>2289533 | 234607<br>4478330 | 488197 | 508293 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 4-Bromofluorobenzene (Surr) | CBZ | Lin1 | +++++<br>384065 | +++++<br>681493 | 78853<br>1195926 | 178351 | 171339 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin1 = Linear 1/conc ISTD |
| Qua = Quadratic ISTD |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57  Calibration End Date: 11/17/2015 06:38  Calibration ID: 31578

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-206828/10 | UX80801.D |
| Level 2 | STD8260 240-206828/9 | UX80800.D |
| Level 3 | STD8260 240-206828/8 | UX80799.D |
| Level 4 | STD8260 240-206828/7 | UX80798.D |
| Level 5 | ICIS 240-206828/6 | UX80797.D |
| Level 6 | STD8260 240-206828/5 | UX80796.D |
| Level 7 | STD8260 240-206828/4 | UX80795.D |
| Level 8 | STD8260 240-206828/3 | UX80794.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | 0.1577<br>0.1509 | 0.1303<br>0.1367 | 0.1661<br>0.1463 | 0.1667 | 0.1431 | Ave | | 0.1497 | | | 0.1000 | 8.8 | | 15.0 | | | |
| Chloromethane | 0.1942<br>0.1878 | 0.1925<br>0.1697 | 0.2201<br>0.1776 | 0.2165 | 0.1804 | Ave | | 0.1924 | | | 0.1000 | 9.3 | | 15.0 | | | |
| Vinyl chloride | 0.1363<br>0.1294 | 0.1352<br>0.1209 | 0.1624<br>0.1282 | 0.1436 | 0.1232 | Ave | | 0.1349 | | | 0.1000 | 9.9 | | 15.0 | | | |
| Butadiene | 0.1536<br>0.1599 | 0.1734<br>0.1504 | 0.2000<br>0.1592 | 0.1773 | 0.1542 | Ave | | 0.1660 | | | | 10.1 | | 15.0 | | | |
| Bromomethane | 0.0711<br>0.0708 | 0.0867<br>0.0508 | 0.0832<br>+++++ | 0.0869 | 0.0689 | Qua | -0.028 | 0.0875 | -0.000362 | | 0.0500 | | | | 0.9980 | | 0.9900 |
| Chloroethane | 0.0904<br>0.0920 | 0.1159<br>0.0792 | 0.1109<br>0.0811 | 0.1096 | 0.0894 | Ave | | 0.0960 | | | | 14.7 | | 15.0 | | | |
| Dichlorofluoromethane | 0.2394<br>0.2116 | 0.2312<br>0.1877 | 0.2630<br>0.1900 | 0.2431 | 0.2122 | Ave | | 0.2223 | | | | 11.9 | | 15.0 | | | |
| Trichlorofluoromethane | 0.2650<br>0.2551 | 0.2457<br>0.2403 | 0.3132<br>0.2226 | 0.2889 | 0.2475 | Ave | | 0.2598 | | | 0.1000 | 11.1 | | 15.0 | | | |
| Ethyl ether | 0.2957<br>0.2391 | 0.2683<br>0.2297 | 0.2653<br>0.2354 | 0.2775 | 0.2368 | Ave | | 0.2560 | | | | 9.4 | | 15.0 | | | |
| Acrolein | 0.0282<br>0.0262 | 0.0290<br>0.0274 | 0.0280<br>0.0276 | 0.0304 | 0.0275 | Ave | | 0.0281 | | | | 4.5 | | 15.0 | | | |
| 1,1-Dichloroethene | 0.2329<br>0.2474 | 0.2133<br>0.2124 | 0.2592<br>0.2131 | 0.2650 | 0.2210 | Ave | | 0.2330 | | | 0.1000 | 9.3 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.1219<br>0.1353 | 0.1071<br>0.1099 | 0.1405<br>0.1205 | 0.1490 | 0.1188 | Ave | | 0.1254 | | | 0.0500 | 11.8 | | 15.0 | | | |
| Acetone | 0.4855<br>0.1448 | 0.3415<br>0.1320 | 0.2016<br>0.1362 | 0.1908 | 0.1481 | Lin1 | 0.7582 | 0.1332 | | | 0.0500 | | | | 0.9980 | | 0.9900 |
| Iodomethane | 0.4693<br>0.4490 | 0.4625<br>0.3967 | 0.5054<br>0.4000 | 0.5032 | 0.4141 | Ave | | 0.4500 | | | | 9.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57 Calibration End Date: 11/17/2015 06:38 Calibration ID: 31578

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon disulfide | 0.5343<br>0.7408 | 0.5526<br>0.6561 | 0.6854<br>0.6642 | 0.7469 | 0.6396 | Ave | | 0.6525 | | | 0.1000 | 11.9 | | 15.0 | | | |
| 3-Chloro-1-propene | 0.0984<br>0.1492 | 0.1171<br>0.1316 | 0.1513<br>0.1372 | 0.1573 | 0.1327 | Ave | | 0.1344 | | | | 14.5 | | 15.0 | | | |
| Methyl acetate | 0.2863<br>0.2624 | 0.2829<br>0.2514 | 0.2782<br>0.2740 | 0.2771 | 0.2467 | Ave | | 0.2699 | | | 0.1000 | 5.4 | | 15.0 | | | |
| Methylene Chloride | +++++<br>0.3009 | 1.1935<br>0.2560 | 0.5815<br>0.2587 | 0.4855 | 0.3682 | Lin1 | 1.9203 | 0.2493 | | | 0.1000 | | | | 0.9970 | | 0.9900 |
| 2-Methyl-2-propanol | 0.0257<br>0.0244 | 0.0221<br>0.0229 | 0.0250<br>0.0196 | 0.0253 | 0.0258 | Ave | | 0.0239 | | | | 9.1 | | 15.0 | | | |
| Acrylonitrile | 0.1247<br>0.1148 | 0.1251<br>0.1071 | 0.1229<br>0.1130 | 0.1230 | 0.1079 | Ave | | 0.1173 | | | | 6.4 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.7152<br>0.7613 | 0.7521<br>0.6900 | 0.7778<br>0.7306 | 0.8069 | 0.6938 | Ave | | 0.7410 | | | 0.1000 | 5.6 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.2599<br>0.2730 | 0.2822<br>0.2397 | 0.3061<br>0.2398 | 0.3127 | 0.2580 | Ave | | 0.2714 | | | 0.1000 | 10.2 | | 15.0 | | | |
| Hexane | 0.0793<br>0.0761 | 0.0674<br>0.0682 | 0.0830<br>0.0701 | 0.0781 | 0.0731 | Ave | | 0.0744 | | | | 7.6 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.5092<br>0.5152 | 0.4879<br>0.4625 | 0.5492<br>0.4800 | 0.5558 | 0.4735 | Ave | | 0.5041 | | | 0.1000 | 6.9 | | 15.0 | | | |
| Vinyl acetate | +++++<br>0.6646 | +++++<br>0.6658 | 0.6233<br>0.7440 | 0.6415 | 0.6274 | Ave | | 0.6611 | | | | 6.7 | | 15.0 | | | |
| 2,2-Dichloropropane | 0.2180<br>0.2975 | 0.2058<br>0.2625 | 0.2605<br>0.2585 | 0.2861 | 0.2463 | Ave | | 0.2544 | | | | 12.2 | | 15.0 | | | |
| 2-Butanone (MEK) | 0.2225<br>0.1793 | 0.1969<br>0.1761 | 0.1845<br>0.1806 | 0.1794 | 0.1772 | Ave | | 0.1871 | | | 0.0500 | 8.4 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.2846<br>0.2988 | 0.2977<br>0.2594 | 0.3287<br>0.2651 | 0.3326 | 0.2731 | Ave | | 0.2925 | | | 0.1000 | 9.4 | | 15.0 | | | |
| Chlorobromomethane | 0.1377<br>0.1505 | 0.1515<br>0.1358 | 0.1647<br>0.1429 | 0.1637 | 0.1401 | Ave | | 0.1484 | | | | 7.6 | | 15.0 | | | |
| Chloroform | 0.4268<br>0.4815 | 0.4597<br>0.4301 | 0.5165<br>0.4447 | 0.5019 | 0.4397 | Ave | | 0.4626 | | | | 7.3 | | 15.0 | | | |
| Tetrahydrofuran | 0.1115<br>0.1050 | 0.1112<br>0.1018 | 0.1081<br>0.1079 | 0.1073 | 0.0974 | Ave | | 0.1063 | | | | 4.5 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.3464<br>0.4190 | 0.3023<br>0.3666 | 0.4016<br>0.3726 | 0.4152 | 0.3653 | Ave | | 0.3736 | | | 0.1000 | 10.4 | | 15.0 | | | |
| Cyclohexane | 0.4788<br>0.5289 | 0.3947<br>0.4551 | 0.5462<br>0.4606 | 0.5452 | 0.4879 | Ave | | 0.4872 | | | 0.1000 | 10.7 | | 15.0 | | | |
| 1,1-Dichloropropene | 0.3313<br>0.3423 | 0.2855<br>0.3043 | 0.3343<br>0.3085 | 0.3478 | 0.3162 | Ave | | 0.3213 | | | | 6.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57 Calibration End Date: 11/17/2015 06:38 Calibration ID: 31578

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon tetrachloride | 0.2800<br>0.3774 | 0.2479<br>0.3373 | 0.3418<br>0.3429 | 0.3557 | 0.3254 | Ave | | 0.3261 | | | 0.1000 | 12.9 | | 15.0 | | | |
| Isobutyl alcohol | 0.0094<br>0.0113 | 0.0096<br>0.0104 | 0.0110<br>0.0097 | 0.0111 | 0.0123 | Ave | | 0.0106 | | | | 9.4 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.4229<br>0.4352 | 0.4216<br>0.3965 | 0.4600<br>0.4193 | 0.4611 | 0.4079 | Ave | | 0.4281 | | | 0.1000 | 5.4 | | 15.0 | | | |
| Benzene | 0.9959<br>0.9946 | 1.0162<br>0.8864 | 1.0849<br>0.9077 | 1.0661 | 0.9252 | Ave | | 0.9846 | | | 0.5000 | 7.4 | | 15.0 | | | |
| n-Heptane | 0.0840<br>0.0830 | 0.0622<br>0.0751 | 0.0874<br>0.0777 | 0.0881 | 0.0761 | Ave | | 0.0792 | | | | 10.7 | | 15.0 | | | |
| Trichloroethene | 0.2820<br>0.3009 | 0.2885<br>0.2752 | 0.3212<br>0.2742 | 0.3179 | 0.2796 | Ave | | 0.2924 | | | 0.2000 | 6.4 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.2491<br>0.2618 | 0.2407<br>0.2337 | 0.2712<br>0.2425 | 0.2650 | 0.2435 | Ave | | 0.2509 | | | 0.1000 | 5.3 | | 15.0 | | | |
| Methylcyclohexane | 0.4420<br>0.4610 | 0.3306<br>0.4074 | 0.4650<br>0.4066 | 0.4787 | 0.4270 | Ave | | 0.4273 | | | 0.1000 | 11.0 | | 15.0 | | | |
| Dibromomethane | 0.1474<br>0.1459 | 0.1413<br>0.1363 | 0.1504<br>0.1429 | 0.1496 | 0.1379 | Ave | | 0.1440 | | | | 3.6 | | 15.0 | | | |
| 1,4-Dioxane | 0.0023<br>0.0021 | 0.0019<br>0.0019 | 0.0021<br>0.0017 | 0.0021 | 0.0020 | Ave | | 0.0020 | | | | 9.7 | | 15.0 | | | |
| Dichlorobromomethane | 0.2635<br>0.3449 | 0.2536<br>0.3214 | 0.3139<br>0.3424 | 0.3150 | 0.3109 | Ave | | 0.3082 | | | 0.1000 | 10.8 | | 15.0 | | | |
| cis-1,3-Dichloropropene | +++++<br>0.3521 | 0.2334<br>0.3415 | 0.2833<br>0.3586 | 0.3080 | 0.3151 | Ave | | 0.3131 | | | 0.1500 | 14.1 | | 15.0 | | | |
| 4-Methyl-2-pentanone (MIBK) | 0.2923<br>0.3312 | 0.2745<br>0.3187 | 0.2978<br>0.3098 | 0.3117 | 0.3346 | Ave | | 0.3088 | | | 0.0500 | 6.5 | | 15.0 | | | |
| Toluene | 1.8864<br>1.5757 | 1.5113<br>1.4192 | 1.6933<br>1.4777 | 1.6655 | 1.4643 | Ave | | 1.5867 | | | 0.4000 | 9.8 | | 15.0 | | | |
| trans-1,3-Dichloropropene | +++++<br>0.4385 | 0.2337<br>0.4379 | 0.3200<br>0.4782 | 0.3409 | 0.3809 | Lin1 | -0.693 | 0.4633 | | | 0.1000 | | | | 0.9970 | | 0.9900 |
| Ethyl methacrylate | +++++<br>0.4136 | 0.2730<br>0.3957 | 0.3479<br>0.4092 | 0.3749 | 0.4119 | Ave | | 0.3752 | | | | 13.6 | | 15.0 | | | |
| 1,1,2-Trichloroethane | 0.3088<br>0.3026 | 0.2744<br>0.2809 | 0.3096<br>0.2932 | 0.3135 | 0.2990 | Ave | | 0.2977 | | | 0.1000 | 4.7 | | 15.0 | | | |
| 1,3-Dichloropropane | 0.5672<br>0.5123 | 0.5253<br>0.4784 | 0.5312<br>0.5032 | 0.5128 | 0.5037 | Ave | | 0.5168 | | | | 5.0 | | 15.0 | | | |
| Tetrachloroethene | 0.3714<br>0.3532 | 0.3279<br>0.3083 | 0.3795<br>0.3130 | 0.3854 | 0.3190 | Ave | | 0.3447 | | | 0.2000 | 9.1 | | 15.0 | | | |
| 2-Hexanone | 0.2824<br>0.3408 | 0.2957<br>0.3306 | 0.3210<br>0.3235 | 0.3289 | 0.3657 | Ave | | 0.3236 | | | 0.0500 | 7.9 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57  Calibration End Date: 11/17/2015 06:38  Calibration ID: 31578

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Chlorodibromomethane | 0.2562<br>0.3664 | 0.2427<br>0.3475 | 0.3196<br>0.3738 | 0.3204 | 0.3304 | Ave | | 0.3196 | | | | 14.9 | | 15.0 | | | |
| Ethylene Dibromide | 0.2737<br>0.3106 | 0.2849<br>0.2923 | 0.2973<br>0.3023 | 0.3115 | 0.3000 | Ave | | 0.2966 | | | | 4.3 | | 15.0 | | | |
| Chlorobenzene | 1.0928<br>1.0194 | 1.0762<br>0.9120 | 1.1237<br>0.9280 | 1.0973 | 0.9949 | Ave | | 1.0305 | | | 0.3000 | 7.8 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.2881<br>0.3973 | 0.3144<br>0.3476 | 0.3689<br>0.3414 | 0.3971 | 0.3655 | Ave | | 0.3525 | | | | 10.8 | | 15.0 | | | |
| Ethylbenzene | 0.5480<br>0.5314 | 0.5133<br>0.4604 | 0.5644<br>0.4559 | 0.5703 | 0.4980 | Ave | | 0.5177 | | | | 8.5 | | 15.0 | | | |
| m-Xylene & p-Xylene | 0.6535<br>0.6442 | 0.6171<br>0.5631 | 0.6729<br>0.5443 | 0.6768 | 0.6077 | Ave | | 0.6225 | | | | 7.9 | | 15.0 | | | |
| o-Xylene | 0.5145<br>0.6240 | 0.5808<br>0.5180 | 0.6833<br>0.4843 | 0.6886 | 0.5985 | Ave | | 0.5865 | | | | 13.2 | | 15.0 | | | |
| Styrene | 0.8513<br>1.0028 | 0.8971<br>0.8774 | 0.9864<br>0.8463 | 1.0235 | 0.9631 | Ave | | 0.9310 | | | 0.3000 | 7.6 | | 15.0 | | | |
| Bromoform | +++++<br>0.1835 | 0.1259<br>0.1774 | 0.1464<br>0.1851 | 0.1617 | 0.1789 | Ave | | 0.1656 | | | 0.1000 | 13.5 | | 15.0 | | | |
| Isopropylbenzene | 1.4084<br>1.6418 | 1.3205<br>1.3639 | 1.6381<br>1.2598 | 1.7376 | 1.5468 | Ave | | 1.4896 | | | 0.1000 | 11.7 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 0.7829<br>0.7849 | 0.7286<br>0.7488 | 0.7834<br>0.7386 | 0.8054 | 0.8681 | Ave | | 0.7801 | | | 0.3000 | 5.7 | | 15.0 | | | |
| Bromobenzene | 1.0340<br>1.0142 | 1.0088<br>0.9466 | 1.0448<br>0.8817 | 1.0236 | 1.0141 | Ave | | 0.9960 | | | | 5.5 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.2579<br>0.2942 | 0.2915<br>0.2784 | 0.3302<br>0.2613 | 0.3065 | 0.3371 | Ave | | 0.2946 | | | | 9.9 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | +++++<br>0.2888 | 0.2087<br>0.3029 | 0.2330<br>0.2952 | 0.2492 | 0.2904 | Ave | | 0.2669 | | | | 13.7 | | 15.0 | | | |
| N-Propylbenzene | 1.0252<br>1.2015 | 1.0431<br>1.0979 | 1.1962<br>0.9826 | 1.2152 | 1.1418 | Ave | | 1.1129 | | | | 8.0 | | 15.0 | | | |
| 2-Chlorotoluene | 0.9156<br>0.9889 | 0.9209<br>0.9041 | 1.0256<br>0.8264 | 1.0538 | 0.9574 | Ave | | 0.9491 | | | | 7.7 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 2.4429<br>3.2490 | 2.5443<br>2.8232 | 3.0645<br>2.6356 | 3.2338 | 3.1315 | Ave | | 2.8906 | | | | 11.1 | | 15.0 | | | |
| 4-Chlorotoluene | 0.9427<br>0.9712 | 0.9558<br>0.8962 | 1.0228<br>0.8315 | 0.9951 | 0.9887 | Ave | | 0.9505 | | | | 6.4 | | 15.0 | | | |
| tert-Butylbenzene | 2.4900<br>3.1690 | 2.3684<br>2.7602 | 2.9068<br>2.8452 | 2.9607 | 3.3023 | Ave | | 2.8503 | | | | 11.0 | | 15.0 | | | |
| 1,2,4-Trimethylbenzene | 2.4595<br>3.1624 | 2.5519<br>2.7522 | 3.1529<br>2.6873 | 3.2452 | 3.0900 | Ave | | 2.8877 | | | | 10.7 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57 Calibration End Date: 11/17/2015 06:38 Calibration ID: 31578

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| sec-Butylbenzene | 3.1227<br>3.9123 | 2.8951<br>3.3639 | 3.6513<br>3.2672 | 3.8300 | 3.7214 | Ave | | 3.4705 | | | | 10.5 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.6842<br>1.6405 | 1.6354<br>1.4695 | 1.7612<br>1.4620 | 1.8153 | 1.6525 | Ave | | 1.6401 | | | 0.6000 | 7.6 | | 15.0 | | | |
| 4-Isopropyltoluene | 2.8485<br>3.4706 | 2.5657<br>2.9553 | 3.2381<br>2.9856 | 3.4588 | 3.3301 | Ave | | 3.1066 | | | | 10.3 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 1.7787<br>1.5984 | 1.7047<br>1.4631 | 1.8304<br>1.4864 | 1.7389 | 1.6851 | Ave | | 1.6607 | | | 0.5000 | 8.0 | | 15.0 | | | |
| n-Butylbenzene | 1.9870<br>2.5464 | 1.8912<br>2.2217 | 2.4002<br>2.3257 | 2.4892 | 2.4004 | Ave | | 2.2827 | | | | 10.3 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.5026<br>1.4792 | 1.4685<br>1.3312 | 1.6088<br>1.4131 | 1.6734 | 1.5236 | Ave | | 1.5000 | | | 0.4000 | 7.1 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | +++++<br>0.1259 | +++++<br>0.1338 | 0.1148<br>0.1600 | 0.1233 | 0.1268 | Ave | | 0.1308 | | | 0.0500 | 11.9 | | 15.0 | | | |
| 1,2,4-Trichlorobenzene | 1.0223<br>1.0300 | 0.9600<br>0.9702 | 1.0532<br>1.0587 | 1.1312 | 0.9906 | Ave | | 1.0270 | | | 0.2000 | 5.4 | | 15.0 | | | |
| Hexachlorobutadiene | 0.4779<br>0.5529 | 0.5049<br>0.4900 | 0.5198<br>0.5536 | 0.5631 | 0.5149 | Ave | | 0.5221 | | | | 6.0 | | 15.0 | | | |
| Naphthalene | 2.0893<br>2.4103 | 2.1684<br>2.3818 | 2.2323<br>2.6445 | 2.4174 | 2.4784 | Ave | | 2.3528 | | | | 7.7 | | 15.0 | | | |
| 1,2,3-Trichlorobenzene | 1.1472<br>0.9798 | 1.0947<br>0.9279 | 1.0807<br>1.0057 | 1.1142 | 1.0122 | Ave | | 1.0453 | | | | 7.2 | | 15.0 | | | |
| Dibromofluoromethane (Surr) | 0.2598<br>0.2446 | 0.2739<br>0.2432 | 0.2740<br>0.2593 | 0.2954 | 0.2586 | Ave | | 0.2636 | | | | 6.5 | | 15.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | 0.3711<br>0.3077 | 0.3491<br>0.3141 | 0.3329<br>0.3371 | 0.3630 | 0.3379 | Ave | | 0.3391 | | | | 6.4 | | 15.0 | | | |
| Toluene-d8 (Surr) | 1.2942<br>1.3135 | 1.4018<br>1.3070 | 1.3980<br>1.3767 | 1.5490 | 1.3645 | Ave | | 1.3756 | | | | 5.9 | | 15.0 | | | |
| 4-Bromofluorobenzene (Surr) | 0.4762<br>0.4031 | 0.4680<br>0.3928 | 0.4664<br>0.3814 | 0.4901 | 0.4453 | Ave | | 0.4404 | | | | 9.5 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57 Calibration End Date: 11/17/2015 06:38 Calibration ID: 31578

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-206828/10 | UX80801.D |
| Level 2 | STD8260 240-206828/9 | UX80800.D |
| Level 3 | STD8260 240-206828/8 | UX80799.D |
| Level 4 | STD8260 240-206828/7 | UX80798.D |
| Level 5 | ICIS 240-206828/6 | UX80797.D |
| Level 6 | STD8260 240-206828/5 | UX80796.D |
| Level 7 | STD8260 240-206828/4 | UX80795.D |
| Level 8 | STD8260 240-206828/3 | UX80794.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 |
| Dichlorodifluoromethane | FB | Ave | 5737<br>280119 | 9191<br>566633 | 29670<br>1079483 | 56206 | 109253 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Chloromethane | FB | Ave | 7065<br>348586 | 13581<br>703686 | 39309<br>1310629 | 73013 | 137749 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Vinyl chloride | FB | Ave | 4959<br>240204 | 9539<br>501084 | 29004<br>946049 | 48421 | 94081 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Butadiene | FB | Ave | 5588<br>296807 | 12231<br>623633 | 35719<br>1174412 | 59769 | 117704 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Bromomethane | FB | Qua | 2588<br>131371 | 6115<br>210597 | 14857<br>+++++ | 29303 | 52623 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>+++++ | 10.0 | 20.0 |
| Chloroethane | FB | Ave | 3288<br>170727 | 8176<br>328192 | 19806<br>598228 | 36948 | 68256 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Dichlorofluoromethane | FB | Ave | 8712<br>392803 | 16311<br>778264 | 46985<br>1401660 | 81963 | 161993 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Trichlorofluoromethane | FB | Ave | 9642<br>473495 | 17329<br>996131 | 55941<br>1642777 | 97391 | 188972 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Ethyl ether | FB | Ave | 10759<br>443795 | 18923<br>952464 | 47396<br>1736824 | 93573 | 180780 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Acrolein | FB | Ave | 5136<br>243414 | 10230<br>567174 | 25028<br>1017791 | 51326 | 105147 | 5.00<br>250 | 10.0<br>500 | 25.0<br>1000 | 50.0 | 100 |
| 1,1-Dichloroethene | FB | Ave | 8474<br>459241 | 15043<br>880676 | 46301<br>1572127 | 89348 | 168747 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroetha ne | FB | Ave | 4434<br>251204 | 7552<br>455619 | 25101<br>889247 | 50223 | 90686 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Acetone | FB | Lin1 | 35330<br>537436 | 48174<br>1094240 | 72029<br>2010360 | 128683 | 226204 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| Iodomethane | FB | Ave | 17077<br>833397 | 32623<br>1644743 | 90282<br>2951678 | 169654 | 316174 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Carbon disulfide | FB | Ave | 19439<br>1375039 | 38977<br>2720161 | 122425<br>4900665 | 251828 | 488300 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57 Calibration End Date: 11/17/2015 06:38 Calibration ID: 31578

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 |
| 3-Chloro-1-propene | FB | Ave | 3581<br>276875 | 8257<br>545754 | 27018<br>1012671 | 53047 | 101348 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Methyl acetate | FB | Ave | 52077<br>2435257 | 99769<br>5212290 | 248439<br>10106918 | 467187 | 941853 | 5.00<br>250 | 10.0<br>500 | 25.0<br>1000 | 50.0 | 100 |
| Methylene Chloride | FB | Lin1 | +++++<br>558441 | 84189<br>1061500 | 103869<br>1909060 | 163680 | 281104 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2-Methyl-2-propanol | FB | Ave | 9366<br>452982 | 15570<br>948826 | 44686<br>1448667 | 85430 | 197240 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100 | 200 |
| Acrylonitrile | FB | Ave | 45380<br>2131066 | 88218<br>4441689 | 219503<br>8338159 | 414751 | 824068 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100 | 200 |
| Methyl tert-butyl ether | FB | Ave | 26024<br>1413028 | 53055<br>2860636 | 138945<br>5390456 | 272063 | 529661 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| trans-1,2-Dichloroethene | FB | Ave | 9455<br>506735 | 19908<br>993744 | 54684<br>1769685 | 105416 | 196980 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Hexane | FB | Ave | 2886<br>141309 | 4752<br>282669 | 14823<br>517517 | 26337 | 55833 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1-Dichloroethane | FB | Ave | 18527<br>956184 | 34414<br>1917420 | 98100<br>3541621 | 187395 | 361480 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Vinyl acetate | FB | Ave | +++++<br>1233486 | +++++<br>2760298 | 111347<br>5489771 | 216302 | 478967 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2,2-Dichloropropane | FB | Ave | 7931<br>552150 | 14517<br>1088257 | 46530<br>1907250 | 96473 | 188004 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2-Butanone (MEK) | FB | Ave | 16194<br>665541 | 27777<br>1460329 | 65921<br>2665491 | 120963 | 270596 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| cis-1,2-Dichloroethene | FB | Ave | 10356<br>554525 | 20997<br>1075316 | 58718<br>1956037 | 112135 | 208539 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Chlorobromomethane | FB | Ave | 5010<br>279400 | 10689<br>562917 | 29412<br>1054199 | 55191 | 106986 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Chloroform | FB | Ave | 15528<br>893689 | 32423<br>1782959 | 92260<br>3281415 | 169220 | 335705 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Tetrahydrofuran | FB | Ave | 8113<br>389703 | 15681<br>843724 | 38619<br>1592365 | 72365 | 148775 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| 1,1,1-Trichloroethane | FB | Ave | 12603<br>777630 | 21326<br>1519876 | 71741<br>2748931 | 140006 | 278923 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Cyclohexane | FB | Ave | 17420<br>981742 | 27843<br>1886623 | 97558<br>3398631 | 183819 | 372493 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1-Dichloropropene | FB | Ave | 12054<br>635276 | 20137<br>1261690 | 59721<br>2276365 | 117251 | 241387 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Carbon tetrachloride | FB | Ave | 10187<br>700474 | 17483<br>1398545 | 61061<br>2530045 | 119946 | 248438 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Isobutyl alcohol | FB | Ave | 8582<br>523630 | 16921<br>1077307 | 49143<br>1782069 | 93390 | 234403 | 25.0<br>1250 | 50.0<br>2500 | 125<br>5000 | 250 | 500 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57 Calibration End Date: 11/17/2015 06:38 Calibration ID: 31578

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 |
| 1,2-Dichloroethane | FB | Ave | 15388<br>807757 | 29737<br>1643955 | 82171<br>3093564 | 155452 | 311422 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Benzene | FB | Ave | 36237<br>1846077 | 71680<br>3674932 | 193794<br>6697613 | 359455 | 706377 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| n-Heptane | FB | Ave | 3056<br>154022 | 4385<br>311272 | 15604<br>573330 | 29704 | 58137 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Trichloroethene | FB | Ave | 10261<br>558435 | 20349<br>1140876 | 57374<br>2022963 | 107182 | 213428 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2-Dichloropropane | FB | Ave | 9064<br>485888 | 16980<br>969024 | 48450<br>1789118 | 89351 | 185929 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Methylcyclohexane | FB | Ave | 16084<br>855704 | 23320<br>1689115 | 83066<br>2999761 | 161394 | 325963 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Dibromomethane | FB | Ave | 5365<br>270786 | 9967<br>565203 | 26864<br>1054082 | 50430 | 105260 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,4-Dioxane | FB | Ave | 1708<br>76930 | 2640<br>158755 | 7383<br>249446 | 14373 | 30291 | 20.0<br>1000 | 40.0<br>2000 | 100<br>4000 | 200 | 400 |
| Dichlorobromomethane | FB | Ave | 9586<br>640159 | 17889<br>1332387 | 56067<br>2526732 | 106203 | 237398 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| cis-1,3-Dichloropropene | FB | Ave | +++++<br>653461 | 16461<br>1415613 | 50612<br>2645605 | 103843 | 240564 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | FB | Ave | 21272<br>1229540 | 38729<br>2642934 | 106406<br>4571653 | 210181 | 510979 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| Toluene | CBZ | Ave | 43668<br>1834046 | 66325<br>3729984 | 188735<br>6673230 | 342571 | 718234 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| trans-1,3-Dichloropropene | CBZ | Lin1 | +++++<br>510340 | 10255<br>1150883 | 35663<br>2159420 | 70111 | 186851 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Ethyl methacrylate | CBZ | Ave | +++++<br>481376 | 11979<br>1040091 | 38775<br>1848148 | 77103 | 202048 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1,2-Trichloroethane | CBZ | Ave | 7148<br>352191 | 12043<br>738242 | 34507<br>1323965 | 64478 | 146642 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,3-Dichloropropane | CBZ | Ave | 13130<br>596292 | 23053<br>1257375 | 59211<br>2272415 | 105470 | 247068 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Tetrachloroethene | CBZ | Ave | 8597<br>411154 | 14390<br>810387 | 42300<br>1413393 | 79267 | 156452 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2-Hexanone | CBZ | Ave | 13076<br>793388 | 25953<br>1737712 | 71545<br>2921908 | 135323 | 358745 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0 | 40.0 |
| Chlorodibromomethane | CBZ | Ave | 5931<br>426433 | 10652<br>913242 | 35627<br>1688291 | 65910 | 162058 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Ethylene Dibromide | CBZ | Ave | 6335<br>361565 | 12502<br>768093 | 33139<br>1365022 | 64070 | 147128 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Chlorobenzene | CBZ | Ave | 25298<br>1186539 | 47229<br>2396834 | 125249<br>4190722 | 225706 | 487975 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57 Calibration End Date: 11/17/2015 06:38 Calibration ID: 31578

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 |
| 1,1,1,2-Tetrachloroethane | CBZ | Ave | 6669<br>462463 | 13798<br>913553 | 41120<br>1541798 | 81683 | 179254 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Ethylbenzene | CBZ | Ave | 12685<br>618509 | 22526<br>1210066 | 62910<br>2059011 | 117300 | 244277 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| m-Xylene & p-Xylene | CBZ | Ave | 15129<br>749855 | 27084<br>1480045 | 74995<br>2457942 | 139205 | 298064 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| o-Xylene | CBZ | Ave | 11910<br>726249 | 25489<br>1361381 | 76156<br>2187083 | 141643 | 293571 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Styrene | CBZ | Ave | 19707<br>1167181 | 39373<br>2306037 | 109946<br>3821988 | 210516 | 472378 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Bromoform | CBZ | Ave | +++++<br>213573 | 5527<br>466222 | 16313<br>836019 | 33263 | 87767 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Isopropylbenzene | CBZ | Ave | 32603<br>1910939 | 57951<br>3584468 | 182575<br>5689483 | 357399 | 758668 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,1,2,2-Tetrachloroethane | DCB | Ave | 6893<br>336344 | 12279<br>662393 | 34383<br>1144797 | 65901 | 159661 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Bromobenzene | DCB | Ave | 9104<br>434624 | 17000<br>837288 | 45858<br>1366657 | 83753 | 186518 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2,3-Trichloropropane | DCB | Ave | 2271<br>126095 | 4912<br>246257 | 14492<br>404973 | 25081 | 62005 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| trans-1,4-Dichloro-2-butene | DCB | Ave | +++++<br>123757 | 3517<br>267916 | 10229<br>457536 | 20387 | 53407 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| N-Propylbenzene | DCB | Ave | 9026<br>514892 | 17578<br>971163 | 52502<br>1523040 | 99423 | 210002 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 2-Chlorotoluene | DCB | Ave | 8061<br>423780 | 15520<br>799743 | 45016<br>1280959 | 86220 | 176088 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,3,5-Trimethylbenzene | DCB | Ave | 21508<br>1392338 | 42878<br>2497328 | 134507<br>4085150 | 264585 | 575963 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 4-Chlorotoluene | DCB | Ave | 8300<br>416184 | 16108<br>792741 | 44893<br>1288832 | 81418 | 181855 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| tert-Butylbenzene | DCB | Ave | 21923<br>1358066 | 39914<br>2441556 | 127585<br>4410077 | 242242 | 607382 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2,4-Trimethylbenzene | DCB | Ave | 21654<br>1355212 | 43005<br>2434463 | 138388<br>4165408 | 265519 | 568337 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| sec-Butylbenzene | DCB | Ave | 27493<br>1676595 | 48789<br>2975559 | 160262<br>5064275 | 313363 | 684456 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,3-Dichlorobenzene | DCB | Ave | 14828<br>703038 | 27560<br>1299865 | 77301<br>2266091 | 148522 | 303942 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 4-Isopropyltoluene | DCB | Ave | 25079<br>1487285 | 43238<br>2614185 | 142128<br>4627675 | 282992 | 612489 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,4-Dichlorobenzene | DCB | Ave | 15660<br>684972 | 28728<br>1294229 | 80342<br>2303873 | 142275 | 309933 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206828

SDG No.:

Instrument ID: A3UX8 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 11/17/2015 03:57 Calibration End Date: 11/17/2015 06:38 Calibration ID: 31578

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4 | LVL 5 |
| n-Butylbenzene | DCB | Ave | 17494<br>1091248 | 31872<br>1965268 | 105348<br>3604810 | 203665 | 441494 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2-Dichlorobenzene | DCB | Ave | 13229<br>633913 | 24747<br>1177499 | 70613<br>2190327 | 136916 | 280228 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | DCB | Ave | +++++<br>53966 | +++++<br>118313 | 5041<br>247973 | 10092 | 23314 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2,4-Trichlorobenzene | DCB | Ave | 9001<br>441403 | 16179<br>858225 | 46228<br>1641074 | 92552 | 182193 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Hexachlorobutadiene | DCB | Ave | 4208<br>236936 | 8508<br>433468 | 22816<br>858018 | 46073 | 94700 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Naphthalene | DCB | Ave | 18395<br>1032934 | 36542<br>2106875 | 97979<br>4099081 | 197789 | 455835 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2,3-Trichlorobenzene | DCB | Ave | 10100<br>419868 | 18448<br>820790 | 47432<br>1558879 | 91164 | 186165 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Dibromofluoromethane (Surr) | FB | Ave | 9452<br>454000 | 19323<br>1008302 | 48938<br>1913539 | 99600 | 197398 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Ave | 13504<br>571219 | 24626<br>1302196 | 59460<br>2487467 | 122377 | 257982 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| Toluene-d8 (Surr) | CBZ | Ave | 29961<br>1528845 | 61519<br>3435042 | 155813<br>6217168 | 318606 | 669279 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |
| 4-Bromofluorobenzene (Surr) | CBZ | Ave | 11024<br>469154 | 20540<br>1032259 | 51987<br>1722622 | 100810 | 218427 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0 | 20.0 |

Curve Type Legend:

Ave = Average ISTD
Lin1 = Linear 1/conc ISTD
Qua = Quadratic ISTD

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-206893/16  Calibration Date: 11/16/2015 14:33

Instrument ID: A3UX10  Calib Start Date: 11/16/2015 12:19

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 11/16/2015 14:10

Lab File ID: UXX3053.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0257 | 0.0243 | | 0.0948 | 0.100 | -5.2 | 20.0 |
| Diisopropyl ether | Ave | 0.2058 | 0.2144 | | 0.0104 | 0.0100 | 4.1 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.3186 | 0.3246 | | 0.0102 | 0.0100 | 1.9 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.6371 | 0.6542 | | 0.0103 | 0.0100 | 2.7 | 20.0 |
| Ethyl acetate | Ave | 0.1936 | 0.1906 | | 0.0197 | 0.0200 | -1.5 | 20.0 |
| Propionitrile | Ave | 0.0318 | 0.0324 | | 0.102 | 0.100 | 1.9 | 20.0 |
| Methacrylonitrile | Ave | 0.1468 | 0.1484 | | 0.101 | 0.100 | 1.0 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.5941 | 0.6019 | | 0.0101 | 0.0100 | 1.3 | 20.0 |
| n-Butanol | Ave | 0.0065 | 0.0057 | | 0.219 | 0.250 | -12.6 | 20.0 |
| Methyl methacrylate | Ave | 0.1691 | 0.1865 | | 0.0221 | 0.0200 | 10.3 | 20.0 |
| 2-Nitropropane | Ave | 0.0384 | 0.0366 | | 0.0190 | 0.0200 | -4.8 | 20.0 |
| 1-Chlorohexane | Ave | 0.3501 | 0.3599 | | 0.0103 | 0.0100 | 2.8 | 20.0 |
| Cyclohexanone | Ave | 0.0129 | 0.0102 | | 0.0789 | 0.100 | -21.1* | 20.0 |
| 1,2,3-Trimethylbenzene | Lin1 | | 2.296 | | 0.00946 | 0.0100 | -5.4 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 0.8458 | 0.8473 | | 0.0100 | 0.0100 | 0.2 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.003 | 0.6222 | | 0.0124 | 0.0200 | -38.0* | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207096/2 Calibration Date: 11/17/2015 10:50

Instrument ID: A3UX10 Calib Start Date: 11/16/2015 09:23

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/16/2015 11:35

Lab File ID: UXX3066.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2491 | 0.2364 | 0.1000 | 0.00949 | 0.0100 | -5.1 | 20.0 |
| Chloromethane | Ave | 0.3073 | 0.2731 | 0.1000 | 0.00889 | 0.0100 | -11.1 | 20.0 |
| Vinyl chloride | Ave | 0.2994 | 0.2876 | 0.1000 | 0.00961 | 0.0100 | -3.9 | 20.0 |
| Butadiene | Ave | 0.3102 | 0.2986 | | 0.00963 | 0.0100 | -3.7 | 20.0 |
| Bromomethane | Ave | 0.2498 | 0.2050 | 0.0500 | 0.00821 | 0.0100 | -17.9 | 20.0 |
| Chloroethane | Ave | 0.2508 | 0.2165 | 0.0500 | 0.00863 | 0.0100 | -13.7 | 20.0 |
| Dichlorofluoromethane | Ave | 0.5829 | 0.4687 | | 0.00804 | 0.0100 | -19.6 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2803 | 0.2754 | 0.1000 | 0.00982 | 0.0100 | -1.8 | 20.0 |
| Ethyl ether | Ave | 0.2326 | 0.2675 | | 0.0115 | 0.0100 | 15.0 | 20.0 |
| Acrolein | Ave | 0.0238 | 0.0218 | | 0.0456 | 0.0500 | -8.7 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2223 | 0.2129 | 0.1000 | 0.00958 | 0.0100 | -4.2 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1658 | 0.1697 | 0.0500 | 0.0102 | 0.0100 | 2.4 | 20.0 |
| Acetone | Lin1 | | 0.0694 | 0.0500 | 0.0212 | 0.0200 | 6.2 | 20.0 |
| Iodomethane | Ave | 0.3742 | 0.3459 | | 0.00924 | 0.0100 | -7.6 | 20.0 |
| Carbon disulfide | Ave | 0.6668 | 0.6195 | 0.1000 | 0.00929 | 0.0100 | -7.1 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1540 | 0.1451 | | 0.00943 | 0.0100 | -5.7 | 20.0 |
| Methyl acetate | Ave | 0.1810 | 0.1705 | 0.1000 | 0.0471 | 0.0500 | -5.8 | 20.0 |
| Methylene Chloride | Lin1 | | 0.2827 | 0.1000 | 0.0109 | 0.0100 | 8.9 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0172 | 0.0145 | | 0.0841 | 0.100 | -15.9 | 20.0 |
| Acrylonitrile | Ave | 0.0902 | 0.0857 | | 0.0949 | 0.100 | -5.1 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7185 | 0.6520 | 0.1000 | 0.00907 | 0.0100 | -9.3 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2610 | 0.2453 | 0.1000 | 0.00940 | 0.0100 | -6.0 | 20.0 |
| Hexane | Ave | 0.0534 | 0.0495 | | 0.00926 | 0.0100 | -7.4 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4582 | 0.4191 | 0.1000 | 0.00915 | 0.0100 | -8.5 | 20.0 |
| Vinyl acetate | Ave | 0.4551 | 0.4093 | | 0.00899 | 0.0100 | -10.1 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2605 | 0.2261 | | 0.00868 | 0.0100 | -13.2 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2789 | 0.2563 | 0.1000 | 0.00919 | 0.0100 | -8.1 | 20.0 |
| 2-Butanone (MEK) | Lin1 | | 0.0917 | 0.0500 | 0.0199 | 0.0200 | -0.6 | 20.0 |
| Bromochloromethane | Ave | 0.1388 | 0.1270 | | 0.00915 | 0.0100 | -8.5 | 20.0 |
| Tetrahydrofuran | Ave | 0.0678 | 0.0617 | | 0.0182 | 0.0200 | -8.9 | 20.0 |
| Chloroform | Ave | 0.4402 | 0.4127 | 0.2000 | 0.00937 | 0.0100 | -6.3 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3307 | 0.3118 | 0.1000 | 0.00943 | 0.0100 | -5.7 | 20.0 |
| Cyclohexane | Ave | 0.3530 | 0.3457 | 0.1000 | 0.00979 | 0.0100 | -2.1 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3118 | 0.2891 | | 0.00927 | 0.0100 | -7.3 | 20.0 |
| Carbon tetrachloride | Ave | 0.2737 | 0.2561 | 0.1000 | 0.00936 | 0.0100 | -6.4 | 20.0 |
| Isobutyl alcohol | Ave | 0.0099 | 0.0086 | | 0.217 | 0.250 | -13.3 | 20.0 |
| Benzene | Ave | 1.051 | 0.9501 | 0.5000 | 0.00904 | 0.0100 | -9.6 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.3543 | 0.3265 | 0.1000 | 0.00922 | 0.0100 | -7.8 | 20.0 |
| n-Heptane | Lin1 | | 0.0393 | | 0.00901 | 0.0100 | -9.9 | 20.0 |
| Trichloroethene | Ave | 0.2661 | 0.2502 | 0.2000 | 0.00940 | 0.0100 | -6.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207096/2 Calibration Date: 11/17/2015 10:50

Instrument ID: A3UX10 Calib Start Date: 11/16/2015 09:23

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/16/2015 11:35

Lab File ID: UXX3066.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.3145 | 0.3071 | 0.1000 | 0.00976 | 0.0100 | -2.4 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2342 | 0.2195 | 0.1000 | 0.00937 | 0.0100 | -6.3 | 20.0 |
| Dibromomethane | Ave | 0.1487 | 0.1399 | | 0.00941 | 0.0100 | -5.9 | 20.0 |
| 1,4-Dioxane | Ave | 0.0025 | 0.0020 | | 0.159 | 0.200 | -20.5* | 20.0 |
| Bromodichloromethane | Ave | 0.2964 | 0.2771 | 0.1000 | 0.00935 | 0.0100 | -6.5 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.1152 | 0.1182 | | 0.0205 | 0.0200 | 2.6 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3280 | 0.3142 | 0.1500 | 0.00958 | 0.0100 | -4.2 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.1898 | 0.1840 | 0.0500 | 0.0194 | 0.0200 | -3.1 | 20.0 |
| Toluene | Ave | 1.490 | 1.379 | 0.4000 | 0.00925 | 0.0100 | -7.5 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3911 | 0.3651 | 0.1000 | 0.00933 | 0.0100 | -6.7 | 20.0 |
| Ethyl methacrylate | Ave | 0.3125 | 0.3268 | | 0.0105 | 0.0100 | 4.6 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2810 | 0.2750 | 0.1000 | 0.00979 | 0.0100 | -2.1 | 20.0 |
| Tetrachloroethene | Ave | 0.2926 | 0.2819 | 0.2000 | 0.00963 | 0.0100 | -3.7 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.5003 | 0.4732 | | 0.00946 | 0.0100 | -5.4 | 20.0 |
| 2-Hexanone | Ave | 0.1698 | 0.1660 | 0.0500 | 0.0196 | 0.0200 | -2.2 | 20.0 |
| Dibromochloromethane | Ave | 0.2817 | 0.2786 | | 0.00989 | 0.0100 | -1.1 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2725 | 0.2697 | | 0.00990 | 0.0100 | -1.0 | 20.0 |
| Chlorobenzene | Ave | 0.9637 | 0.8860 | 0.3000 | 0.00919 | 0.0100 | -8.1 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3193 | 0.3052 | | 0.00956 | 0.0100 | -4.4 | 20.0 |
| Ethylbenzene | Ave | 0.5026 | 0.4674 | | 0.00930 | 0.0100 | -7.0 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6012 | 0.5753 | | 0.00957 | 0.0100 | -4.3 | 20.0 |
| o-Xylene | Ave | 0.5969 | 0.5727 | | 0.00959 | 0.0100 | -4.1 | 20.0 |
| Styrene | Ave | 0.9732 | 0.9609 | 0.3000 | 0.00987 | 0.0100 | -1.3 | 20.0 |
| Bromoform | Ave | 0.1822 | 0.1685 | 0.1000 | 0.00925 | 0.0100 | -7.5 | 20.0 |
| Isopropylbenzene | Ave | 1.510 | 1.416 | 0.1000 | 0.00938 | 0.0100 | -6.2 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.6201 | 0.6202 | 0.3000 | 0.0100 | 0.0100 | 0.0 | 20.0 |
| Bromobenzene | Ave | 0.7341 | 0.7196 | | 0.00980 | 0.0100 | -2.0 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.2180 | 0.2136 | | 0.00980 | 0.0100 | -2.0 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.1903 | 0.1809 | | 0.00951 | 0.0100 | -4.9 | 20.0 |
| N-Propylbenzene | Ave | 0.7036 | 0.7266 | | 0.0103 | 0.0100 | 3.3 | 20.0 |
| 2-Chlorotoluene | Ave | 0.6570 | 0.6320 | | 0.00962 | 0.0100 | -3.8 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 2.032 | 2.049 | | 0.0101 | 0.0100 | 0.8 | 20.0 |
| 4-Chlorotoluene | Ave | 0.6937 | 0.6980 | | 0.0101 | 0.0100 | 0.6 | 20.0 |
| tert-Butylbenzene | Ave | 1.781 | 1.732 | | 0.00973 | 0.0100 | -2.7 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 2.223 | 2.138 | | 0.00962 | 0.0100 | -3.8 | 20.0 |
| sec-Butylbenzene | Ave | 2.317 | 2.247 | | 0.00970 | 0.0100 | -3.0 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.375 | 1.279 | 0.6000 | 0.00930 | 0.0100 | -7.0 | 20.0 |
| p-Isopropyltoluene | Ave | 2.071 | 1.988 | | 0.00960 | 0.0100 | -4.0 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.447 | 1.327 | 0.5000 | 0.00917 | 0.0100 | -8.3 | 20.0 |
| n-Butylbenzene | Ave | 1.594 | 1.471 | | 0.00923 | 0.0100 | -7.7 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.329 | 1.207 | 0.4000 | 0.00908 | 0.0100 | -9.2 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207096/2  Calibration Date: 11/17/2015 10:50

Instrument ID: A3UX10  Calib Start Date: 11/16/2015 09:23

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 11/16/2015 11:35

Lab File ID: UXX3066.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.1084 | 0.0841 | 0.0500 | 0.00776 | 0.0100 | -22.4* | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.8230 | 0.5017 | 0.2000 | 0.00610 | 0.0100 | -39.0* | 20.0 |
| Hexachlorobutadiene | Ave | 0.2912 | 0.1969 | | 0.00676 | 0.0100 | -32.4* | 20.0 |
| Naphthalene | Ave | 2.023 | 1.184 | | 0.00585 | 0.0100 | -41.5* | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.7870 | 0.4455 | | 0.00566 | 0.0100 | -43.4* | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2341 | 0.1968 | | 0.0103 | 0.0123 | -15.9 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2958 | 0.2462 | | 0.0102 | 0.0123 | -16.8 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.265 | 1.134 | | 0.0110 | 0.0123 | -10.4 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 0.4298 | 0.3549 | | 0.0102 | 0.0123 | -17.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207096/3  Calibration Date: 11/17/2015 11:11

Instrument ID: A3UX10  Calib Start Date: 11/16/2015 12:19

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 11/16/2015 14:10

Lab File ID: UXX3067.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0257 | 0.0201 | | 0.0782 | 0.100 | -21.8* | 20.0 |
| Diisopropyl ether | Ave | 0.2058 | 0.1952 | | 0.00948 | 0.0100 | -5.2 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.3186 | 0.3024 | | 0.00949 | 0.0100 | -5.1 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.6371 | 0.5951 | | 0.00934 | 0.0100 | -6.6 | 20.0 |
| Ethyl acetate | Ave | 0.1936 | 0.1740 | | 0.0180 | 0.0200 | -10.1 | 20.0 |
| Propionitrile | Ave | 0.0318 | 0.0283 | | 0.0890 | 0.100 | -11.0 | 20.0 |
| Methacrylonitrile | Ave | 0.1468 | 0.1365 | | 0.0930 | 0.100 | -7.0 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.5941 | 0.5644 | | 0.00950 | 0.0100 | -5.0 | 20.0 |
| n-Butanol | Ave | 0.0065 | 0.0050 | | 0.191 | 0.250 | -23.7* | 20.0 |
| Methyl methacrylate | Ave | 0.1691 | 0.1692 | | 0.0200 | 0.0200 | 0.0 | 20.0 |
| 2-Nitropropane | Ave | 0.0384 | 0.0390 | | 0.0203 | 0.0200 | 1.3 | 20.0 |
| 1-Chlorohexane | Ave | 0.3501 | 0.3321 | | 0.00948 | 0.0100 | -5.2 | 20.0 |
| Cyclohexanone | Ave | 0.0129 | 0.0094 | | 0.0725 | 0.100 | -27.5* | 20.0 |
| 1,2,3-Trimethylbenzene | Lin1 | | 2.096 | | 0.00868 | 0.0100 | -13.2 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 0.8458 | 0.7547 | | 0.00892 | 0.0100 | -10.8 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.003 | 0.5145 | | 0.0103 | 0.0200 | -48.7* | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-196026/13 Calibration Date: 09/02/2015 18:07

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 144945.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2422 | 0.2902 | 0.1000 | 0.0240 | 0.0200 | 19.8 | 20.0 |
| Chloromethane | Ave | 0.3787 | 0.4264 | 0.1000 | 0.0225 | 0.0200 | 12.6 | 20.0 |
| Vinyl chloride | Ave | 0.2747 | 0.2991 | 0.1000 | 0.0218 | 0.0200 | 8.9 | 20.0 |
| Butadiene | Ave | 0.3003 | 0.3280 | | 0.0218 | 0.0200 | 9.2 | 20.0 |
| Bromomethane | Ave | 0.1383 | 0.1509 | 0.0500 | 0.0218 | 0.0200 | 9.1 | 20.0 |
| Chloroethane | Ave | 0.1448 | 0.1548 | 0.0500 | 0.0214 | 0.0200 | 6.9 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3245 | 0.3543 | | 0.0218 | 0.0200 | 9.2 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2852 | 0.3259 | 0.1000 | 0.0229 | 0.0200 | 14.3 | 20.0 |
| Ethyl ether | Ave | 0.1804 | 0.2164 | | 0.0240 | 0.0200 | 19.9 | 20.0 |
| Acrolein | Ave | 0.0180 | 0.0235 | | 0.130 | 0.100 | 30.4* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2217 | 0.2411 | 0.1000 | 0.0218 | 0.0200 | 8.8 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1625 | 0.1744 | 0.0500 | 0.0215 | 0.0200 | 7.4 | 20.0 |
| Acetone | Lin1 | | 0.0867 | 0.0500 | 0.0453 | 0.0400 | 13.2 | 20.0 |
| Iodomethane | Ave | 0.3194 | 0.3613 | | 0.0226 | 0.0200 | 13.1 | 20.0 |
| Carbon disulfide | Ave | 0.6714 | 0.6482 | 0.1000 | 0.0193 | 0.0200 | -3.4 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1430 | 0.1394 | | 0.0195 | 0.0200 | -2.5 | 20.0 |
| Methyl acetate | Ave | 0.1567 | 0.1835 | 0.1000 | 0.117 | 0.100 | 17.1 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3235 | 0.1000 | 0.0217 | 0.0200 | 8.6 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0198 | 0.0221 | | 0.223 | 0.200 | 11.3 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2515 | 0.2789 | 0.1000 | 0.0222 | 0.0200 | 10.9 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.5650 | 0.6276 | 0.1000 | 0.0222 | 0.0200 | 11.1 | 20.0 |
| Acrylonitrile | Ave | 0.0757 | 0.0862 | | 0.228 | 0.200 | 13.8 | 20.0 |
| Hexane | Ave | 0.0816 | 0.0814 | | 0.0200 | 0.0200 | -0.1 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4925 | 0.5567 | 0.1000 | 0.0226 | 0.0200 | 13.0 | 20.0 |
| Vinyl acetate | Lin1 | | 0.0217 | | 0.0149 | 0.0200 | -25.6* | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2107 | 0.2260 | | 0.0214 | 0.0200 | 7.2 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2654 | 0.2924 | 0.1000 | 0.0220 | 0.0200 | 10.2 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0962 | 0.1043 | 0.0500 | 0.0433 | 0.0400 | 8.4 | 20.0 |
| Bromochloromethane | Ave | 0.0967 | 0.1063 | | 0.0220 | 0.0200 | 9.9 | 20.0 |
| Tetrahydrofuran | Ave | 0.0665 | 0.0705 | | 0.0424 | 0.0400 | 6.0 | 20.0 |
| Chloroform | Ave | 0.3979 | 0.4367 | 0.2000 | 0.0219 | 0.0200 | 9.7 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3127 | 0.3465 | 0.1000 | 0.0222 | 0.0200 | 10.8 | 20.0 |
| Cyclohexane | Ave | 0.5900 | 0.6590 | 0.1000 | 0.0223 | 0.0200 | 11.7 | 20.0 |
| Carbon tetrachloride | Ave | 0.2541 | 0.2901 | 0.1000 | 0.0228 | 0.0200 | 14.2 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3331 | 0.3884 | | 0.0233 | 0.0200 | 16.6 | 20.0 |
| Benzene | Ave | 0.999 | 1.113 | 0.5000 | 0.0223 | 0.0200 | 11.4 | 20.0 |
| Isobutyl alcohol | Ave | 0.0071 | 0.0082 | | 0.573 | 0.500 | 14.6 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.2792 | 0.3100 | 0.1000 | 0.0222 | 0.0200 | 11.0 | 20.0 |
| n-Heptane | Ave | 0.0682 | 0.0773 | | 0.0227 | 0.0200 | 13.4 | 20.0 |
| Trichloroethene | Ave | 0.2336 | 0.2694 | 0.2000 | 0.0231 | 0.0200 | 15.3 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-196026/13 Calibration Date: 09/02/2015 18:07

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 144945.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.4826 | 0.5370 | 0.1000 | 0.0223 | 0.0200 | 11.3 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2493 | 0.2803 | 0.1000 | 0.0225 | 0.0200 | 12.5 | 20.0 |
| Dibromomethane | Ave | 0.1040 | 0.1205 | | 0.0232 | 0.0200 | 15.9 | 20.0 |
| 1,4-Dioxane | Ave | 0.0017 | 0.0019 | | 0.462 | 0.400 | 15.4 | 20.0 |
| Bromodichloromethane | Ave | 0.2508 | 0.2588 | 0.1000 | 0.0206 | 0.0200 | 3.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0933 | 0.1063 | | 0.0228 | 0.0200 | 14.0 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3090 | 0.3296 | 0.1500 | 0.0213 | 0.0200 | 6.7 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3102 | 0.3490 | 0.0500 | 0.0450 | 0.0400 | 12.5 | 20.0 |
| Toluene | Ave | 1.617 | 1.748 | 0.4000 | 0.0216 | 0.0200 | 8.1 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3680 | 0.4131 | 0.1000 | 0.0224 | 0.0200 | 12.2 | 20.0 |
| Ethyl methacrylate | Ave | 0.3012 | 0.3245 | | 0.0215 | 0.0200 | 7.7 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2333 | 0.2665 | 0.1000 | 0.0228 | 0.0200 | 14.2 | 20.0 |
| Tetrachloroethene | Ave | 0.2833 | 0.3198 | 0.2000 | 0.0226 | 0.0200 | 12.9 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4264 | 0.4668 | | 0.0219 | 0.0200 | 9.5 | 20.0 |
| 2-Hexanone | Ave | 0.2007 | 0.2198 | 0.0500 | 0.0438 | 0.0400 | 9.5 | 20.0 |
| Dibromochloromethane | Lin1 | | 0.2290 | | 0.0193 | 0.0200 | -3.6 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2138 | 0.2523 | | 0.0236 | 0.0200 | 18.0 | 20.0 |
| Chlorobenzene | Ave | 0.9298 | 1.009 | 0.3000 | 0.0217 | 0.0200 | 8.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.2902 | 0.3344 | | 0.0231 | 0.0200 | 15.3 | 20.0 |
| Ethylbenzene | Ave | 0.5296 | 0.5798 | | 0.0219 | 0.0200 | 9.5 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6468 | 0.6849 | | 0.0212 | 0.0200 | 5.9 | 20.0 |
| o-Xylene | Ave | 0.6346 | 0.7209 | | 0.0227 | 0.0200 | 13.6 | 20.0 |
| Styrene | Ave | 0.9556 | 0.9735 | 0.3000 | 0.0204 | 0.0200 | 1.9 | 20.0 |
| Bromoform | Lin1 | | 0.1061 | 0.1000 | 0.0173 | 0.0200 | -13.4 | 20.0 |
| Isopropylbenzene | Ave | 1.742 | 1.946 | 0.1000 | 0.0223 | 0.0200 | 11.7 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.5958 | 0.6826 | 0.3000 | 0.0229 | 0.0200 | 14.6 | 20.0 |
| Bromobenzene | Ave | 0.7042 | 0.7858 | | 0.0223 | 0.0200 | 11.6 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1763 | 0.2024 | | 0.0230 | 0.0200 | 14.8 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2152 | 0.2320 | | 0.0216 | 0.0200 | 7.8 | 20.0 |
| N-Propylbenzene | Ave | 0.9769 | 1.052 | | 0.0215 | 0.0200 | 7.7 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8014 | 0.8578 | | 0.0214 | 0.0200 | 7.0 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.102 | 3.417 | | 0.0220 | 0.0200 | 10.2 | 20.0 |
| 4-Chlorotoluene | Ave | 0.7776 | 0.8277 | | 0.0213 | 0.0200 | 6.5 | 20.0 |
| tert-Butylbenzene | Ave | 2.752 | 3.074 | | 0.0223 | 0.0200 | 11.7 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.081 | 3.372 | | 0.0219 | 0.0200 | 9.5 | 20.0 |
| sec-Butylbenzene | Ave | 4.194 | 4.751 | | 0.0227 | 0.0200 | 13.3 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.474 | 1.583 | 0.6000 | 0.0215 | 0.0200 | 7.4 | 20.0 |
| p-Isopropyltoluene | Ave | 3.242 | 3.585 | | 0.0221 | 0.0200 | 10.6 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.484 | 1.542 | 0.5000 | 0.0208 | 0.0200 | 3.9 | 20.0 |
| n-Butylbenzene | Ave | 3.148 | 3.388 | | 0.0215 | 0.0200 | 7.6 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.421 | 1.521 | 0.4000 | 0.0214 | 0.0200 | 7.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-196026/13 Calibration Date: 09/02/2015 18:07

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 144945.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0871 | 0.0879 | 0.0500 | 0.0202 | 0.0200 | 0.9 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.9733 | 0.9859 | 0.2000 | 0.0203 | 0.0200 | 1.3 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5396 | 0.6222 | | 0.0231 | 0.0200 | 15.3 | 20.0 |
| Naphthalene | Ave | 2.263 | 2.304 | | 0.0204 | 0.0200 | 1.8 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.9223 | 0.9287 | | 0.0201 | 0.0200 | 0.7 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2154 | 0.2054 | | 0.0477 | 0.0500 | -4.6 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2516 | 0.2368 | | 0.0471 | 0.0500 | -5.9 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.539 | 1.437 | | 0.0467 | 0.0500 | -6.6 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.147 | 1.045 | | 0.0455 | 0.0500 | -8.9 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207343/3 Calibration Date: 11/18/2015 11:15

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 146532.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2422 | 0.2743 | 0.1000 | 0.0226 | 0.0200 | 13.2 | 20.0 |
| Chloromethane | Ave | 0.3787 | 0.3080 | 0.1000 | 0.0163 | 0.0200 | -18.7 | 20.0 |
| Vinyl chloride | Ave | 0.2747 | 0.2356 | 0.1000 | 0.0172 | 0.0200 | -14.2 | 20.0 |
| Butadiene | Ave | 0.3003 | 0.2712 | | 0.0181 | 0.0200 | -9.7 | 20.0 |
| Bromomethane | Ave | 0.1383 | 0.1411 | 0.0500 | 0.0204 | 0.0200 | 2.0 | 20.0 |
| Chloroethane | Ave | 0.1448 | 0.1451 | 0.0500 | 0.0200 | 0.0200 | 0.2 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3245 | 0.3506 | | 0.0216 | 0.0200 | 8.0 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2852 | 0.3577 | 0.1000 | 0.0251 | 0.0200 | 25.4* | 20.0 |
| Ethyl ether | Ave | 0.1804 | 0.1589 | | 0.0440 | 0.0500 | -11.9 | 20.0 |
| Acrolein | Ave | 0.0180 | 0.0126 | | 0.175 | 0.250 | -30.2* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2217 | 0.2165 | 0.1000 | 0.0488 | 0.0500 | -2.3 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1625 | 0.1927 | 0.0500 | 0.0593 | 0.0500 | 18.6 | 20.0 |
| Acetone | Lin1 | | 0.0618 | 0.0500 | 0.0886 | 0.100 | -11.4 | 20.0 |
| Iodomethane | Ave | 0.3194 | 0.3376 | | 0.0529 | 0.0500 | 5.7 | 20.0 |
| Carbon disulfide | Ave | 0.6714 | 0.6767 | 0.1000 | 0.0504 | 0.0500 | 0.8 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1430 | 0.1353 | | 0.0473 | 0.0500 | -5.4 | 20.0 |
| Methyl acetate | Ave | 0.1567 | 0.1205 | 0.1000 | 0.192 | 0.250 | -23.1* | 20.0 |
| Methylene Chloride | Lin1 | | 0.2477 | 0.1000 | 0.0479 | 0.0500 | -4.3 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0198 | 0.0136 | | 0.342 | 0.500 | -31.5* | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2515 | 0.2649 | 0.1000 | 0.0527 | 0.0500 | 5.3 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.5650 | 0.5222 | 0.1000 | 0.0462 | 0.0500 | -7.6 | 20.0 |
| Acrylonitrile | Ave | 0.0757 | 0.0603 | | 0.398 | 0.500 | -20.4* | 20.0 |
| Hexane | Ave | 0.0816 | 0.0749 | | 0.0459 | 0.0500 | -8.2 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4925 | 0.4761 | 0.1000 | 0.0483 | 0.0500 | -3.3 | 20.0 |
| Vinyl acetate | Lin1 | | 0.0303 | | 0.0505 | 0.0500 | 1.0 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2107 | 0.2751 | | 0.0653 | 0.0500 | 30.5* | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2654 | 0.2785 | 0.1000 | 0.0525 | 0.0500 | 5.0 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0962 | 0.0740 | 0.0500 | 0.0769 | 0.100 | -23.1* | 20.0 |
| Bromochloromethane | Ave | 0.0967 | 0.1104 | | 0.0571 | 0.0500 | 14.2 | 20.0 |
| Tetrahydrofuran | Ave | 0.0665 | 0.0462 | | 0.0695 | 0.100 | -30.5* | 20.0 |
| Chloroform | Ave | 0.3979 | 0.4355 | 0.2000 | 0.0547 | 0.0500 | 9.4 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3127 | 0.4000 | 0.1000 | 0.0639 | 0.0500 | 27.9* | 20.0 |
| Cyclohexane | Ave | 0.5900 | 0.5412 | 0.1000 | 0.0459 | 0.0500 | -8.3 | 20.0 |
| Carbon tetrachloride | Ave | 0.2541 | 0.3285 | 0.1000 | 0.0647 | 0.0500 | 29.3* | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3331 | 0.3388 | | 0.0508 | 0.0500 | 1.7 | 20.0 |
| Benzene | Ave | 0.999 | 0.9835 | 0.5000 | 0.0492 | 0.0500 | -1.5 | 20.0 |
| Isobutyl alcohol | Ave | 0.0071 | 0.0052 | | 0.908 | 1.25 | -27.3* | 20.0 |
| 1,2-Dichloroethane | Ave | 0.2792 | 0.2889 | 0.1000 | 0.0517 | 0.0500 | 3.5 | 20.0 |
| n-Heptane | Ave | 0.0682 | 0.0727 | | 0.0533 | 0.0500 | 6.7 | 20.0 |
| Trichloroethene | Ave | 0.2336 | 0.2604 | 0.2000 | 0.0557 | 0.0500 | 11.5 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207343/3  Calibration Date: 11/18/2015 11:15

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146532.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.4826 | 0.5023 | 0.1000 | 0.0520 | 0.0500 | 4.1 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2493 | 0.2278 | 0.1000 | 0.0457 | 0.0500 | -8.6 | 20.0 |
| Dibromomethane | Ave | 0.1040 | 0.1094 | | 0.0526 | 0.0500 | 5.2 | 20.0 |
| 1,4-Dioxane | Ave | 0.0017 | 0.0014 | | 0.836 | 1.00 | -16.4 | 20.0 |
| Bromodichloromethane | Ave | 0.2508 | 0.2760 | 0.1000 | 0.0550 | 0.0500 | 10.0 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0933 | 0.0698 | | 0.0748 | 0.100 | -25.2* | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3090 | 0.3009 | 0.1500 | 0.0487 | 0.0500 | -2.6 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3102 | 0.2335 | 0.0500 | 0.0753 | 0.100 | -24.7* | 20.0 |
| Toluene | Ave | 1.617 | 1.694 | 0.4000 | 0.0524 | 0.0500 | 4.8 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3680 | 0.3630 | 0.1000 | 0.0493 | 0.0500 | -1.4 | 20.0 |
| Ethyl methacrylate | Ave | 0.3012 | 0.2672 | | 0.0444 | 0.0500 | -11.3 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2333 | 0.2490 | 0.1000 | 0.0533 | 0.0500 | 6.7 | 20.0 |
| Tetrachloroethene | Ave | 0.2833 | 0.3070 | 0.2000 | 0.0542 | 0.0500 | 8.4 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4264 | 0.3980 | | 0.0467 | 0.0500 | -6.7 | 20.0 |
| 2-Hexanone | Ave | 0.2007 | 0.1527 | 0.0500 | 0.0761 | 0.100 | -23.9* | 20.0 |
| Dibromochloromethane | Lin1 | | 0.2555 | | 0.0515 | 0.0500 | 3.0 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2138 | 0.2247 | | 0.0526 | 0.0500 | 5.1 | 20.0 |
| Chlorobenzene | Ave | 0.9298 | 0.9529 | 0.3000 | 0.0512 | 0.0500 | 2.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.2902 | 0.3328 | | 0.0573 | 0.0500 | 14.7 | 20.0 |
| Ethylbenzene | Ave | 0.5296 | 0.5477 | | 0.0517 | 0.0500 | 3.4 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6468 | 0.6760 | | 0.0523 | 0.0500 | 4.5 | 20.0 |
| o-Xylene | Ave | 0.6346 | 0.6629 | | 0.0522 | 0.0500 | 4.5 | 20.0 |
| Styrene | Ave | 0.9556 | 1.008 | 0.3000 | 0.0528 | 0.0500 | 5.5 | 20.0 |
| Bromoform | Lin1 | | 0.1226 | 0.1000 | 0.0481 | 0.0500 | -3.8 | 20.0 |
| Isopropylbenzene | Ave | 1.742 | 1.875 | 0.1000 | 0.0538 | 0.0500 | 7.7 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.5958 | 0.5638 | 0.3000 | 0.0473 | 0.0500 | -5.4 | 20.0 |
| Bromobenzene | Ave | 0.7042 | 0.7788 | | 0.0553 | 0.0500 | 10.6 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1763 | 0.1750 | | 0.0496 | 0.0500 | -0.8 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2152 | 0.1802 | | 0.0419 | 0.0500 | -16.3 | 20.0 |
| N-Propylbenzene | Ave | 0.9769 | 1.043 | | 0.0534 | 0.0500 | 6.8 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8014 | 0.8396 | | 0.0524 | 0.0500 | 4.8 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.102 | 3.299 | | 0.0532 | 0.0500 | 6.4 | 20.0 |
| 4-Chlorotoluene | Ave | 0.7776 | 0.8246 | | 0.0530 | 0.0500 | 6.1 | 20.0 |
| tert-Butylbenzene | Ave | 2.752 | 2.986 | | 0.0543 | 0.0500 | 8.5 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.081 | 3.265 | | 0.0530 | 0.0500 | 6.0 | 20.0 |
| sec-Butylbenzene | Ave | 4.194 | 4.313 | | 0.0514 | 0.0500 | 2.8 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.474 | 1.446 | 0.6000 | 0.0490 | 0.0500 | -1.9 | 20.0 |
| p-Isopropyltoluene | Ave | 3.242 | 3.567 | | 0.0550 | 0.0500 | 10.0 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.484 | 1.409 | 0.5000 | 0.0475 | 0.0500 | -5.1 | 20.0 |
| n-Butylbenzene | Ave | 3.148 | 3.291 | | 0.0523 | 0.0500 | 4.6 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.421 | 1.375 | 0.4000 | 0.0484 | 0.0500 | -3.2 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207343/3  Calibration Date: 11/18/2015 11:15

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146532.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0871 | 0.0837 | 0.0500 | 0.0480 | 0.0500 | -3.9 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.9733 | 0.9553 | 0.2000 | 0.0491 | 0.0500 | -1.8 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5396 | 0.6017 | | 0.0558 | 0.0500 | 11.5 | 20.0 |
| Naphthalene | Ave | 2.263 | 1.911 | | 0.0422 | 0.0500 | -15.6 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.9223 | 0.8900 | | 0.0482 | 0.0500 | -3.5 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2154 | 0.2266 | | 0.0526 | 0.0500 | 5.2 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2516 | 0.2573 | | 0.0511 | 0.0500 | 2.3 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.539 | 1.528 | | 0.0497 | 0.0500 | -0.7 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.147 | 1.042 | | 0.0454 | 0.0500 | -9.2 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207343/4  Calibration Date: 11/18/2015 11:37

Instrument ID: A3UX14  Calib Start Date: 08/13/2015 17:50

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 08/13/2015 20:42

Lab File ID: 146533.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0224 | 0.0193 | | 0.431 | 0.500 | -13.7 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.4341 | 0.4232 | | 0.0487 | 0.0500 | -2.5 | 20.0 |
| Diisopropyl ether | Ave | 0.2079 | 0.1901 | | 0.0457 | 0.0500 | -8.6 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8262 | 0.6633 | | 0.0401 | 0.0500 | -19.7 | 20.0 |
| Ethyl acetate | Ave | 0.1606 | 0.1204 | | 0.0750 | 0.100 | -25.0* | 20.0 |
| Propionitrile | Ave | 0.0278 | 0.0200 | | 0.360 | 0.500 | -27.9* | 20.0 |
| Methacrylonitrile | Ave | 0.1070 | 0.0950 | | 0.444 | 0.500 | -11.3 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.7020 | 0.5399 | | 0.0385 | 0.0500 | -23.1* | 20.0 |
| n-Butanol | Ave | 0.0105 | 0.0059 | | 0.701 | 1.25 | -43.9* | 20.0 |
| Methyl methacrylate | Ave | 0.1432 | 0.1197 | | 0.0836 | 0.100 | -16.4 | 20.0 |
| 2-Nitropropane | Ave | 0.0224 | 0.0209 | | 0.0932 | 0.100 | -6.8 | 20.0 |
| 1-Chlorohexane | Ave | 0.4504 | 0.4262 | | 0.0473 | 0.0500 | -5.4 | 20.0 |
| Cyclohexanone | Ave | 0.0123 | 0.0067 | | 0.271 | 0.500 | -45.7* | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.034 | 3.030 | | 0.0499 | 0.0500 | -0.1 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.096 | 1.094 | | 0.0499 | 0.0500 | -0.2 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.220 | 1.179 | | 0.0966 | 0.100 | -3.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207560/3 Calibration Date: 11/19/2015 11:25

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 146565.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2422 | 0.2594 | 0.1000 | 0.0214 | 0.0200 | 7.1 | 20.0 |
| Chloromethane | Ave | 0.3787 | 0.2950 | 0.1000 | 0.0156 | 0.0200 | -22.1* | 20.0 |
| Vinyl chloride | Ave | 0.2747 | 0.2254 | 0.1000 | 0.0164 | 0.0200 | -17.9 | 20.0 |
| Butadiene | Ave | 0.3003 | 0.2536 | | 0.0169 | 0.0200 | -15.6 | 20.0 |
| Bromomethane | Ave | 0.1383 | 0.1318 | 0.0500 | 0.0191 | 0.0200 | -4.7 | 20.0 |
| Chloroethane | Ave | 0.1448 | 0.1347 | 0.0500 | 0.0186 | 0.0200 | -6.9 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3245 | 0.3178 | | 0.0196 | 0.0200 | -2.1 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2852 | 0.3283 | 0.1000 | 0.0230 | 0.0200 | 15.1 | 20.0 |
| Ethyl ether | Ave | 0.1804 | 0.1510 | | 0.0419 | 0.0500 | -16.3 | 20.0 |
| Acrolein | Ave | 0.0180 | 0.0117 | | 0.162 | 0.250 | -35.2* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2217 | 0.2175 | 0.1000 | 0.0491 | 0.0500 | -1.9 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1625 | 0.1828 | 0.0500 | 0.0563 | 0.0500 | 12.5 | 20.0 |
| Acetone | Lin1 | | 0.0551 | 0.0500 | 0.0778 | 0.100 | -22.2* | 20.0 |
| Iodomethane | Ave | 0.3194 | 0.3091 | | 0.0484 | 0.0500 | -3.2 | 20.0 |
| Carbon disulfide | Ave | 0.6714 | 0.5932 | 0.1000 | 0.0442 | 0.0500 | -11.6 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1430 | 0.1276 | | 0.0446 | 0.0500 | -10.8 | 20.0 |
| Methyl acetate | Ave | 0.1567 | 0.1152 | 0.1000 | 0.184 | 0.250 | -26.5* | 20.0 |
| Methylene Chloride | Lin1 | | 0.2254 | 0.1000 | 0.0429 | 0.0500 | -14.1 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0198 | 0.0144 | | 0.363 | 0.500 | -27.4* | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2515 | 0.2472 | 0.1000 | 0.0491 | 0.0500 | -1.7 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.5650 | 0.4965 | 0.1000 | 0.0439 | 0.0500 | -12.1 | 20.0 |
| Acrylonitrile | Ave | 0.0757 | 0.0576 | | 0.380 | 0.500 | -23.9* | 20.0 |
| Hexane | Ave | 0.0816 | 0.0656 | | 0.0402 | 0.0500 | -19.5 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4925 | 0.4319 | 0.1000 | 0.0438 | 0.0500 | -12.3 | 20.0 |
| Vinyl acetate | Lin1 | | 0.0296 | | 0.0493 | 0.0500 | -1.4 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2107 | 0.2392 | | 0.0567 | 0.0500 | 13.5 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2654 | 0.2596 | 0.1000 | 0.0489 | 0.0500 | -2.2 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0962 | 0.0688 | 0.0500 | 0.0715 | 0.100 | -28.5* | 20.0 |
| Bromochloromethane | Ave | 0.0967 | 0.1028 | | 0.0531 | 0.0500 | 6.2 | 20.0 |
| Tetrahydrofuran | Ave | 0.0665 | 0.0465 | | 0.0699 | 0.100 | -30.1* | 20.0 |
| Chloroform | Ave | 0.3979 | 0.4100 | 0.2000 | 0.0515 | 0.0500 | 3.0 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3127 | 0.3621 | 0.1000 | 0.0579 | 0.0500 | 15.8 | 20.0 |
| Cyclohexane | Ave | 0.5900 | 0.5046 | 0.1000 | 0.0428 | 0.0500 | -14.5 | 20.0 |
| Carbon tetrachloride | Ave | 0.2541 | 0.2999 | 0.1000 | 0.0590 | 0.0500 | 18.1 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3331 | 0.3169 | | 0.0476 | 0.0500 | -4.9 | 20.0 |
| Benzene | Ave | 0.999 | 0.9115 | 0.5000 | 0.0456 | 0.0500 | -8.8 | 20.0 |
| Isobutyl alcohol | Ave | 0.0071 | 0.0057 | | 0.990 | 1.25 | -20.8* | 20.0 |
| 1,2-Dichloroethane | Ave | 0.2792 | 0.2734 | 0.1000 | 0.0490 | 0.0500 | -2.1 | 20.0 |
| n-Heptane | Ave | 0.0682 | 0.0681 | | 0.0499 | 0.0500 | -0.1 | 20.0 |
| Trichloroethene | Ave | 0.2336 | 0.2375 | 0.2000 | 0.0508 | 0.0500 | 1.7 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207560/3  Calibration Date: 11/19/2015 11:25

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146565.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.4826 | 0.4586 | 0.1000 | 0.0475 | 0.0500 | -5.0 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2493 | 0.2168 | 0.1000 | 0.0435 | 0.0500 | -13.0 | 20.0 |
| Dibromomethane | Ave | 0.1040 | 0.1010 | | 0.0486 | 0.0500 | -2.9 | 20.0 |
| 1,4-Dioxane | Ave | 0.0017 | 0.0014 | | 0.820 | 1.00 | -18.0 | 20.0 |
| Bromodichloromethane | Ave | 0.2508 | 0.2588 | 0.1000 | 0.0516 | 0.0500 | 3.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0933 | 0.0668 | | 0.0716 | 0.100 | -28.4* | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3090 | 0.2831 | 0.1500 | 0.0458 | 0.0500 | -8.4 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3102 | 0.2257 | 0.0500 | 0.0727 | 0.100 | -27.3* | 20.0 |
| Toluene | Ave | 1.617 | 1.530 | 0.4000 | 0.0473 | 0.0500 | -5.4 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3680 | 0.3350 | 0.1000 | 0.0455 | 0.0500 | -9.0 | 20.0 |
| Ethyl methacrylate | Ave | 0.3012 | 0.2581 | | 0.0429 | 0.0500 | -14.3 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2333 | 0.2272 | 0.1000 | 0.0487 | 0.0500 | -2.6 | 20.0 |
| Tetrachloroethene | Ave | 0.2833 | 0.2778 | 0.2000 | 0.0490 | 0.0500 | -1.9 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4264 | 0.3771 | | 0.0442 | 0.0500 | -11.6 | 20.0 |
| 2-Hexanone | Ave | 0.2007 | 0.1456 | 0.0500 | 0.0726 | 0.100 | -27.4* | 20.0 |
| Dibromochloromethane | Lin1 | | 0.2388 | | 0.0482 | 0.0500 | -3.6 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2138 | 0.2096 | | 0.0490 | 0.0500 | -2.0 | 20.0 |
| Chlorobenzene | Ave | 0.9298 | 0.8780 | 0.3000 | 0.0472 | 0.0500 | -5.6 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.2902 | 0.2962 | | 0.0510 | 0.0500 | 2.1 | 20.0 |
| Ethylbenzene | Ave | 0.5296 | 0.4808 | | 0.0454 | 0.0500 | -9.2 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6468 | 0.6095 | | 0.0471 | 0.0500 | -5.8 | 20.0 |
| o-Xylene | Ave | 0.6346 | 0.6111 | | 0.0481 | 0.0500 | -3.7 | 20.0 |
| Styrene | Ave | 0.9556 | 0.9219 | 0.3000 | 0.0482 | 0.0500 | -3.5 | 20.0 |
| Bromoform | Lin1 | | 0.1184 | 0.1000 | 0.0465 | 0.0500 | -7.0 | 20.0 |
| Isopropylbenzene | Ave | 1.742 | 1.716 | 0.1000 | 0.0493 | 0.0500 | -1.4 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.5958 | 0.5463 | 0.3000 | 0.0458 | 0.0500 | -8.3 | 20.0 |
| Bromobenzene | Ave | 0.7042 | 0.7072 | | 0.0502 | 0.0500 | 0.4 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1763 | 0.1657 | | 0.0470 | 0.0500 | -6.0 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2152 | 0.1745 | | 0.0405 | 0.0500 | -18.9 | 20.0 |
| N-Propylbenzene | Ave | 0.9769 | 0.9698 | | 0.0496 | 0.0500 | -0.7 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8014 | 0.7736 | | 0.0483 | 0.0500 | -3.5 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.102 | 3.074 | | 0.0495 | 0.0500 | -0.9 | 20.0 |
| 4-Chlorotoluene | Ave | 0.7776 | 0.7509 | | 0.0483 | 0.0500 | -3.4 | 20.0 |
| tert-Butylbenzene | Ave | 2.752 | 2.719 | | 0.0494 | 0.0500 | -1.2 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.081 | 3.050 | | 0.0495 | 0.0500 | -1.0 | 20.0 |
| sec-Butylbenzene | Ave | 4.194 | 4.041 | | 0.0482 | 0.0500 | -3.7 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.474 | 1.350 | 0.6000 | 0.0458 | 0.0500 | -8.4 | 20.0 |
| p-Isopropyltoluene | Ave | 3.242 | 3.307 | | 0.0510 | 0.0500 | 2.0 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.484 | 1.309 | 0.5000 | 0.0441 | 0.0500 | -11.8 | 20.0 |
| n-Butylbenzene | Ave | 3.148 | 3.111 | | 0.0494 | 0.0500 | -1.2 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.421 | 1.277 | 0.4000 | 0.0449 | 0.0500 | -10.2 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207560/3  Calibration Date: 11/19/2015 11:25

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146565.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0871 | 0.0923 | 0.0500 | 0.0530 | 0.0500 | 5.9 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.9733 | 0.9496 | 0.2000 | 0.0488 | 0.0500 | -2.4 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5396 | 0.5693 | | 0.0528 | 0.0500 | 5.5 | 20.0 |
| Naphthalene | Ave | 2.263 | 1.959 | | 0.0433 | 0.0500 | -13.4 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.9223 | 0.8779 | | 0.0476 | 0.0500 | -4.8 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2154 | 0.2144 | | 0.0498 | 0.0500 | -0.4 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2516 | 0.2456 | | 0.0488 | 0.0500 | -2.4 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.539 | 1.397 | | 0.0454 | 0.0500 | -9.2 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.147 | 0.9335 | | 0.0407 | 0.0500 | -18.7 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207560/4  Calibration Date: 11/19/2015 11:46

Instrument ID: A3UX14  Calib Start Date: 08/13/2015 17:50

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 08/13/2015 20:42

Lab File ID: 146566.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0224 | 0.0212 | | 0.474 | 0.500 | -5.2 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.4341 | 0.4465 | | 0.0514 | 0.0500 | 2.9 | 20.0 |
| Diisopropyl ether | Ave | 0.2079 | 0.2047 | | 0.0492 | 0.0500 | -1.6 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8262 | 0.7266 | | 0.0440 | 0.0500 | -12.1 | 20.0 |
| Ethyl acetate | Ave | 0.1606 | 0.1435 | | 0.0893 | 0.100 | -10.7 | 20.0 |
| Propionitrile | Ave | 0.0278 | 0.0225 | | 0.405 | 0.500 | -19.0 | 20.0 |
| Methacrylonitrile | Ave | 0.1070 | 0.1050 | | 0.491 | 0.500 | -1.8 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.7020 | 0.5897 | | 0.0420 | 0.0500 | -16.0 | 20.0 |
| n-Butanol | Ave | 0.0105 | 0.0068 | | 0.812 | 1.25 | -35.0* | 20.0 |
| Methyl methacrylate | Ave | 0.1432 | 0.1432 | | 0.100 | 0.100 | -0.0 | 20.0 |
| 2-Nitropropane | Ave | 0.0224 | 0.0246 | | 0.110 | 0.100 | 9.5 | 20.0 |
| 1-Chlorohexane | Ave | 0.4504 | 0.4898 | | 0.0544 | 0.0500 | 8.8 | 20.0 |
| Cyclohexanone | Ave | 0.0123 | 0.0081 | | 0.329 | 0.500 | -34.1* | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.034 | 3.185 | | 0.0525 | 0.0500 | 5.0 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.096 | 1.216 | | 0.0555 | 0.0500 | 11.0 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.220 | 1.325 | | 0.109 | 0.100 | 8.7 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-206828/11 Calibration Date: 11/17/2015 06:59

Instrument ID: A3UX8 Calib Start Date: 07/30/2015 02:35

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 07/30/2015 05:04

Lab File ID: UX80802.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 2-Chloroethyl vinyl ether | Lin1 | 0.0079 | | | 0.00120 | 0.0250 | -100.0* | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-206828/11 Calibration Date: 11/17/2015 06:59

Instrument ID: A3UX8 Calib Start Date: 11/17/2015 03:57

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/17/2015 06:38

Lab File ID: UX80802.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.1497 | 0.1487 | 0.1000 | 0.0248 | 0.0250 | -0.7 | 20.0 |
| Chloromethane | Ave | 0.1924 | 0.1902 | 0.1000 | 0.0247 | 0.0250 | -1.1 | 20.0 |
| Vinyl chloride | Ave | 0.1349 | 0.1294 | 0.1000 | 0.0240 | 0.0250 | -4.1 | 20.0 |
| Butadiene | Ave | 0.1660 | 0.1604 | | 0.0242 | 0.0250 | -3.4 | 20.0 |
| Bromomethane | Qua | | 0.0753 | 0.0500 | 0.0243 | 0.0250 | -2.8 | 20.0 |
| Chloroethane | Ave | 0.0960 | 0.0898 | 0.0500 | 0.0234 | 0.0250 | -6.5 | 20.0 |
| Dichlorofluoromethane | Ave | 0.2223 | 0.2152 | | 0.0242 | 0.0250 | -3.2 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2598 | 0.2482 | 0.1000 | 0.0239 | 0.0250 | -4.5 | 20.0 |
| Ethyl ether | Ave | 0.2560 | 0.2382 | | 0.0233 | 0.0250 | -7.0 | 20.0 |
| Acrolein | Ave | 0.0281 | 0.0481 | | 0.214 | 0.125 | 71.5* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2330 | 0.2192 | 0.1000 | 0.0235 | 0.0250 | -5.9 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1254 | 0.1210 | 0.0500 | 0.0241 | 0.0250 | -3.5 | 20.0 |
| Acetone | Lin1 | | 0.1530 | 0.0500 | 0.0517 | 0.0500 | 3.5 | 20.0 |
| Iodomethane | Ave | 0.4500 | 0.4532 | | 0.0252 | 0.0250 | 0.7 | 20.0 |
| Carbon disulfide | Ave | 0.6525 | 0.6398 | 0.1000 | 0.0245 | 0.0250 | -1.9 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1344 | 0.1267 | | 0.0236 | 0.0250 | -5.7 | 20.0 |
| Methyl acetate | Ave | 0.2699 | 0.2755 | 0.1000 | 0.128 | 0.125 | 2.1 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3265 | 0.1000 | 0.0250 | 0.0250 | 0.2 | 20.0 |
| 2-Methyl-2-propanol | Ave | 0.0239 | 0.0253 | | 0.265 | 0.250 | 6.1 | 20.0 |
| Acrylonitrile | Ave | 0.1173 | 0.1201 | | 0.256 | 0.250 | 2.3 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7410 | 0.7723 | 0.1000 | 0.0261 | 0.0250 | 4.2 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2714 | 0.2760 | 0.1000 | 0.0254 | 0.0250 | 1.7 | 20.0 |
| Hexane | Ave | 0.0744 | 0.0700 | | 0.0235 | 0.0250 | -6.0 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.5041 | 0.5104 | 0.1000 | 0.0253 | 0.0250 | 1.2 | 20.0 |
| Vinyl acetate | Ave | 0.6611 | 0.5554 | | 0.0210 | 0.0250 | -16.0 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2544 | 0.2361 | | 0.0232 | 0.0250 | -7.2 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.1871 | 0.1796 | 0.0500 | 0.0480 | 0.0500 | -4.0 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2925 | 0.2942 | 0.1000 | 0.0251 | 0.0250 | 0.6 | 20.0 |
| Chlorobromomethane | Ave | 0.1484 | 0.1583 | | 0.0267 | 0.0250 | 6.7 | 20.0 |
| Chloroform | Ave | 0.4626 | 0.4752 | 0.2000 | 0.0257 | 0.0250 | 2.7 | 20.0 |
| Tetrahydrofuran | Ave | 0.1063 | 0.1086 | | 0.0511 | 0.0500 | 2.2 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3736 | 0.3733 | 0.1000 | 0.0250 | 0.0250 | -0.0 | 20.0 |
| Cyclohexane | Ave | 0.4872 | 0.4831 | 0.1000 | 0.0248 | 0.0250 | -0.8 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3213 | 0.3342 | | 0.0260 | 0.0250 | 4.0 | 20.0 |
| Carbon tetrachloride | Ave | 0.3261 | 0.3293 | 0.1000 | 0.0253 | 0.0250 | 1.0 | 20.0 |
| Isobutyl alcohol | Ave | 0.0106 | 0.0120 | | 0.708 | 0.625 | 13.3 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.4281 | 0.4336 | 0.1000 | 0.0253 | 0.0250 | 1.3 | 20.0 |
| Benzene | Ave | 0.9846 | 0.9662 | 0.5000 | 0.0245 | 0.0250 | -1.9 | 20.0 |
| n-Heptane | Ave | 0.0792 | 0.0760 | | 0.0240 | 0.0250 | -4.0 | 20.0 |
| Trichloroethene | Ave | 0.2924 | 0.2992 | 0.2000 | 0.0256 | 0.0250 | 2.3 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-206828/11 Calibration Date: 11/17/2015 06:59

Instrument ID: A3UX8 Calib Start Date: 11/17/2015 03:57

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/17/2015 06:38

Lab File ID: UX80802.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dichloropropane | Ave | 0.2509 | 0.2584 | 0.1000 | 0.0257 | 0.0250 | 3.0 | 20.0 |
| Methylcyclohexane | Ave | 0.4273 | 0.4210 | 0.1000 | 0.0246 | 0.0250 | -1.5 | 20.0 |
| 1,4-Dioxane | Ave | 0.0020 | 0.0022 | | 0.540 | 0.500 | 7.9 | 20.0 |
| Dibromomethane | Ave | 0.1440 | 0.1535 | | 0.0267 | 0.0250 | 6.7 | 20.0 |
| Dichlorobromomethane | Ave | 0.3082 | 0.3152 | 0.1000 | 0.0256 | 0.0250 | 2.3 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3131 | 0.3178 | 0.1500 | 0.0254 | 0.0250 | 1.5 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3088 | 0.3277 | 0.0500 | 0.0530 | 0.0500 | 6.1 | 20.0 |
| Toluene | Ave | 1.587 | 1.548 | 0.4000 | 0.0244 | 0.0250 | -2.5 | 20.0 |
| trans-1,3-Dichloropropene | Lin1 | | 0.4040 | 0.1000 | 0.0233 | 0.0250 | -6.8 | 20.0 |
| Ethyl methacrylate | Ave | 0.3752 | 0.4000 | | 0.0267 | 0.0250 | 6.6 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2977 | 0.3062 | 0.1000 | 0.0257 | 0.0250 | 2.8 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.5168 | 0.4991 | | 0.0241 | 0.0250 | -3.4 | 20.0 |
| Tetrachloroethene | Ave | 0.3447 | 0.3302 | 0.2000 | 0.0240 | 0.0250 | -4.2 | 20.0 |
| 2-Hexanone | Ave | 0.3236 | 0.3267 | 0.0500 | 0.0505 | 0.0500 | 1.0 | 20.0 |
| Chlorodibromomethane | Ave | 0.3196 | 0.3345 | | 0.0262 | 0.0250 | 4.7 | 20.0 |
| Ethylene Dibromide | Ave | 0.2966 | 0.3056 | | 0.0258 | 0.0250 | 3.0 | 20.0 |
| Chlorobenzene | Ave | 1.031 | 1.004 | 0.3000 | 0.0244 | 0.0250 | -2.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3525 | 0.3753 | | 0.0266 | 0.0250 | 6.5 | 20.0 |
| Ethylbenzene | Ave | 0.5177 | 0.5036 | | 0.0243 | 0.0250 | -2.7 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6225 | 0.6141 | | 0.0247 | 0.0250 | -1.3 | 20.0 |
| o-Xylene | Ave | 0.5865 | 0.6119 | | 0.0261 | 0.0250 | 4.3 | 20.0 |
| Styrene | Ave | 0.9310 | 0.9171 | 0.3000 | 0.0246 | 0.0250 | -1.5 | 20.0 |
| Bromoform | Ave | 0.1656 | 0.1683 | 0.1000 | 0.0254 | 0.0250 | 1.6 | 20.0 |
| Isopropylbenzene | Ave | 1.490 | 1.580 | 0.1000 | 0.0265 | 0.0250 | 6.1 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.7801 | 0.8109 | 0.3000 | 0.0260 | 0.0250 | 3.9 | 20.0 |
| Bromobenzene | Ave | 0.996 | 0.9830 | | 0.0247 | 0.0250 | -1.3 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.2946 | 0.3172 | | 0.0269 | 0.0250 | 7.7 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2669 | 0.2469 | | 0.0231 | 0.0250 | -7.5 | 20.0 |
| N-Propylbenzene | Ave | 1.113 | 1.110 | | 0.0249 | 0.0250 | -0.3 | 20.0 |
| 2-Chlorotoluene | Ave | 0.9491 | 0.9748 | | 0.0257 | 0.0250 | 2.7 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 2.891 | 3.052 | | 0.0264 | 0.0250 | 5.6 | 20.0 |
| 4-Chlorotoluene | Ave | 0.9505 | 0.9426 | | 0.0248 | 0.0250 | -0.8 | 20.0 |
| tert-Butylbenzene | Ave | 2.850 | 3.215 | | 0.0282 | 0.0250 | 12.8 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 2.888 | 2.991 | | 0.0259 | 0.0250 | 3.6 | 20.0 |
| sec-Butylbenzene | Ave | 3.470 | 3.656 | | 0.0263 | 0.0250 | 5.4 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.640 | 1.554 | 0.6000 | 0.0237 | 0.0250 | -5.2 | 20.0 |
| 4-Isopropyltoluene | Ave | 3.107 | 3.217 | | 0.0259 | 0.0250 | 3.6 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.661 | 1.574 | 0.5000 | 0.0237 | 0.0250 | -5.2 | 20.0 |
| n-Butylbenzene | Ave | 2.283 | 2.279 | | 0.0250 | 0.0250 | -0.2 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.500 | 1.482 | 0.4000 | 0.0247 | 0.0250 | -1.2 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | Ave | 0.1308 | 0.1156 | 0.0500 | 0.0221 | 0.0250 | -11.6 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-206828/11  Calibration Date: 11/17/2015 06:59

Instrument ID: A3UX8  Calib Start Date: 11/17/2015 03:57

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 11/17/2015 06:38

Lab File ID: UX80802.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2,4-Trichlorobenzene | Ave | 1.027 | 0.9519 | 0.2000 | 0.0232 | 0.0250 | -7.3 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5221 | 0.5029 | | 0.0241 | 0.0250 | -3.7 | 20.0 |
| Naphthalene | Ave | 2.353 | 2.321 | | 0.0247 | 0.0250 | -1.4 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 1.045 | 0.9599 | | 0.0230 | 0.0250 | -8.2 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2636 | 0.2556 | | 0.0485 | 0.0500 | -3.0 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.3391 | 0.3242 | | 0.0478 | 0.0500 | -4.4 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.376 | 1.301 | | 0.0473 | 0.0500 | -5.4 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 0.4404 | 0.4113 | | 0.0467 | 0.0500 | -6.6 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207720/4  Calibration Date: 11/20/2015 03:43

Instrument ID: A3UX8  Calib Start Date: 07/30/2015 02:35

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 07/30/2015 05:04

Lab File ID: UX80875.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 2-Chloroethyl vinyl ether | Lin1 | 0.0079 | 0.0038 | | | 0.0400 | | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207720/4 Calibration Date: 11/20/2015 03:43

Instrument ID: A3UX8 Calib Start Date: 11/17/2015 03:57

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/17/2015 06:38

Lab File ID: UX80875.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.1497 | 0.1863 | 0.1000 | 0.0249 | 0.0200 | 24.4* | 20.0 |
| Chloromethane | Ave | 0.1924 | 0.2006 | 0.1000 | 0.0209 | 0.0200 | 4.3 | 20.0 |
| Vinyl chloride | Ave | 0.1349 | 0.1429 | 0.1000 | 0.0212 | 0.0200 | 5.9 | 20.0 |
| Butadiene | Ave | 0.1660 | 0.1756 | | 0.0212 | 0.0200 | 5.8 | 20.0 |
| Bromomethane | Qua | | 0.0795 | 0.0500 | 0.0202 | 0.0200 | 1.0 | 20.0 |
| Chloroethane | Ave | 0.0960 | 0.0976 | 0.0500 | 0.0203 | 0.0200 | 1.7 | 20.0 |
| Dichlorofluoromethane | Ave | 0.2223 | 0.2305 | | 0.0207 | 0.0200 | 3.7 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2598 | 0.2786 | 0.1000 | 0.0214 | 0.0200 | 7.2 | 20.0 |
| Ethyl ether | Ave | 0.2560 | 0.2179 | | 0.0170 | 0.0200 | -14.9 | 20.0 |
| Acrolein | Ave | 0.0281 | 0.0232 | | 0.0825 | 0.100 | -17.5 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2330 | 0.2331 | 0.1000 | 0.0200 | 0.0200 | 0.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1254 | 0.1315 | 0.0500 | 0.0210 | 0.0200 | 4.9 | 20.0 |
| Acetone | Lin1 | | 0.1565 | 0.0500 | 0.0413 | 0.0400 | 3.2 | 20.0 |
| Iodomethane | Ave | 0.4500 | 0.4287 | | 0.0191 | 0.0200 | -4.7 | 20.0 |
| Carbon disulfide | Ave | 0.6525 | 0.6573 | 0.1000 | 0.0201 | 0.0200 | 0.7 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1344 | 0.1277 | | 0.0190 | 0.0200 | -4.9 | 20.0 |
| Methyl acetate | Ave | 0.2699 | 0.2467 | 0.1000 | 0.0914 | 0.100 | -8.6 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3185 | 0.1000 | 0.0178 | 0.0200 | -10.8 | 20.0 |
| 2-Methyl-2-propanol | Ave | 0.0239 | 0.0217 | | 0.182 | 0.200 | -9.2 | 20.0 |
| Acrylonitrile | Ave | 0.1173 | 0.1146 | | 0.195 | 0.200 | -2.3 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7410 | 0.6535 | 0.1000 | 0.0176 | 0.0200 | -11.8 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2714 | 0.2632 | 0.1000 | 0.0194 | 0.0200 | -3.0 | 20.0 |
| Hexane | Ave | 0.0744 | 0.0662 | | 0.0178 | 0.0200 | -11.0 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.5041 | 0.4671 | 0.1000 | 0.0185 | 0.0200 | -7.4 | 20.0 |
| Vinyl acetate | Ave | 0.6611 | 0.5455 | | 0.0165 | 0.0200 | -17.5 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.1871 | 0.1882 | 0.0500 | 0.0402 | 0.0400 | 0.6 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2544 | 0.2047 | | 0.0161 | 0.0200 | -19.6 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2925 | 0.2814 | 0.1000 | 0.0192 | 0.0200 | -3.8 | 20.0 |
| Chlorobromomethane | Ave | 0.1484 | 0.1465 | | 0.0198 | 0.0200 | -1.2 | 20.0 |
| Chloroform | Ave | 0.4626 | 0.4536 | 0.2000 | 0.0196 | 0.0200 | -2.0 | 20.0 |
| Tetrahydrofuran | Ave | 0.1063 | 0.1036 | | 0.0390 | 0.0400 | -2.5 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3736 | 0.3557 | 0.1000 | 0.0190 | 0.0200 | -4.8 | 20.0 |
| Cyclohexane | Ave | 0.4872 | 0.4985 | 0.1000 | 0.0205 | 0.0200 | 2.3 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3213 | 0.3081 | | 0.0192 | 0.0200 | -4.1 | 20.0 |
| Carbon tetrachloride | Ave | 0.3261 | 0.3144 | 0.1000 | 0.0193 | 0.0200 | -3.6 | 20.0 |
| Isobutyl alcohol | Ave | 0.0106 | 0.0108 | | 0.512 | 0.500 | 2.4 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.4281 | 0.4189 | 0.1000 | 0.0196 | 0.0200 | -2.1 | 20.0 |
| Benzene | Ave | 0.9846 | 0.9081 | 0.5000 | 0.0184 | 0.0200 | -7.8 | 20.0 |
| n-Heptane | Ave | 0.0792 | 0.0710 | | 0.0179 | 0.0200 | -10.4 | 20.0 |
| Trichloroethene | Ave | 0.2924 | 0.2827 | 0.2000 | 0.0193 | 0.0200 | -3.3 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207720/4  Calibration Date: 11/20/2015 03:43

Instrument ID: A3UX8  Calib Start Date: 11/17/2015 03:57

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 11/17/2015 06:38

Lab File ID: UX80875.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dichloropropane | Ave | 0.2509 | 0.2367 | 0.1000 | 0.0189 | 0.0200 | -5.7 | 20.0 |
| Methylcyclohexane | Ave | 0.4273 | 0.4299 | 0.1000 | 0.0201 | 0.0200 | 0.6 | 20.0 |
| 1,4-Dioxane | Ave | 0.0020 | 0.0023 | | 0.456 | 0.400 | 13.9 | 20.0 |
| Dibromomethane | Ave | 0.1440 | 0.1486 | | 0.0206 | 0.0200 | 3.2 | 20.0 |
| Dichlorobromomethane | Ave | 0.3082 | 0.3205 | 0.1000 | 0.0208 | 0.0200 | 4.0 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3131 | 0.2724 | 0.1500 | 0.0174 | 0.0200 | -13.0 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3088 | 0.3266 | 0.0500 | 0.0423 | 0.0400 | 5.8 | 20.0 |
| Toluene | Ave | 1.587 | 1.524 | 0.4000 | 0.0192 | 0.0200 | -3.9 | 20.0 |
| trans-1,3-Dichloropropene | Lin1 | | 0.3191 | 0.1000 | 0.0153 | 0.0200 | -23.6* | 20.0 |
| Ethyl methacrylate | Ave | 0.3752 | 0.3593 | | 0.0192 | 0.0200 | -4.2 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2977 | 0.3075 | 0.1000 | 0.0207 | 0.0200 | 3.3 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.5168 | 0.5032 | | 0.0195 | 0.0200 | -2.6 | 20.0 |
| Tetrachloroethene | Ave | 0.3447 | 0.3327 | 0.2000 | 0.0193 | 0.0200 | -3.5 | 20.0 |
| 2-Hexanone | Ave | 0.3236 | 0.3359 | 0.0500 | 0.0415 | 0.0400 | 3.8 | 20.0 |
| Chlorodibromomethane | Ave | 0.3196 | 0.3464 | | 0.0217 | 0.0200 | 8.4 | 20.0 |
| Ethylene Dibromide | Ave | 0.2966 | 0.2984 | | 0.0201 | 0.0200 | 0.6 | 20.0 |
| Chlorobenzene | Ave | 1.031 | 0.998 | 0.3000 | 0.0194 | 0.0200 | -3.2 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3525 | 0.3718 | | 0.0211 | 0.0200 | 5.5 | 20.0 |
| Ethylbenzene | Ave | 0.5177 | 0.5161 | | 0.0199 | 0.0200 | -0.3 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6225 | 0.6189 | | 0.0199 | 0.0200 | -0.6 | 20.0 |
| o-Xylene | Ave | 0.5865 | 0.6208 | | 0.0212 | 0.0200 | 5.9 | 20.0 |
| Styrene | Ave | 0.9310 | 0.9611 | 0.3000 | 0.0206 | 0.0200 | 3.2 | 20.0 |
| Bromoform | Ave | 0.1656 | 0.1683 | 0.1000 | 0.0203 | 0.0200 | 1.7 | 20.0 |
| Isopropylbenzene | Ave | 1.490 | 1.623 | 0.1000 | 0.0218 | 0.0200 | 9.0 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.7801 | 0.8674 | 0.3000 | 0.0222 | 0.0200 | 11.2 | 20.0 |
| Bromobenzene | Ave | 0.996 | 1.000 | | 0.0201 | 0.0200 | 0.4 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.2946 | 0.3326 | | 0.0226 | 0.0200 | 12.9 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2669 | 0.2611 | | 0.0196 | 0.0200 | -2.2 | 20.0 |
| N-Propylbenzene | Ave | 1.113 | 1.191 | | 0.0214 | 0.0200 | 7.0 | 20.0 |
| 2-Chlorotoluene | Ave | 0.9491 | 1.023 | | 0.0215 | 0.0200 | 7.7 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 2.891 | 3.219 | | 0.0223 | 0.0200 | 11.4 | 20.0 |
| 4-Chlorotoluene | Ave | 0.9505 | 0.9870 | | 0.0208 | 0.0200 | 3.8 | 20.0 |
| tert-Butylbenzene | Ave | 2.850 | 3.110 | | 0.0218 | 0.0200 | 9.1 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 2.888 | 3.208 | | 0.0222 | 0.0200 | 11.1 | 20.0 |
| sec-Butylbenzene | Ave | 3.470 | 3.824 | | 0.0220 | 0.0200 | 10.2 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.640 | 1.713 | 0.6000 | 0.0209 | 0.0200 | 4.4 | 20.0 |
| 4-Isopropyltoluene | Ave | 3.107 | 3.404 | | 0.0219 | 0.0200 | 9.6 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.661 | 1.676 | 0.5000 | 0.0202 | 0.0200 | 0.9 | 20.0 |
| n-Butylbenzene | Ave | 2.283 | 2.453 | | 0.0215 | 0.0200 | 7.4 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.500 | 1.594 | 0.4000 | 0.0213 | 0.0200 | 6.3 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | Ave | 0.1308 | 0.1167 | 0.0500 | 0.0178 | 0.0200 | -10.8 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207720/4  Calibration Date: 11/20/2015 03:43

Instrument ID: A3UX8  Calib Start Date: 11/17/2015 03:57

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 11/17/2015 06:38

Lab File ID: UX80875.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2,4-Trichlorobenzene | Ave | 1.027 | 1.050 | 0.2000 | 0.0204 | 0.0200 | 2.2 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5221 | 0.5513 | | 0.0211 | 0.0200 | 5.6 | 20.0 |
| Naphthalene | Ave | 2.353 | 2.463 | | 0.0209 | 0.0200 | 4.7 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 1.045 | 1.056 | | 0.0202 | 0.0200 | 1.0 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2636 | 0.2631 | | 0.0499 | 0.0500 | -0.2 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.3391 | 0.3119 | | 0.0460 | 0.0500 | -8.0 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.376 | 1.371 | | 0.0498 | 0.0500 | -0.3 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 0.4404 | 0.4124 | | 0.0468 | 0.0500 | -6.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207720/5  Calibration Date: 11/20/2015 04:04

Instrument ID: A3UX8  Calib Start Date: 07/15/2015 01:51

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 07/15/2015 03:38

Lab File ID: UX80876.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0523 | 0.0357 | | 0.136 | 0.200 | -31.9* | 20.0 |
| Isopropyl ether | Ave | 0.2112 | 0.1991 | | 0.0188 | 0.0200 | -5.8 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.5368 | 0.4273 | | 0.0159 | 0.0200 | -20.4* | 20.0 |
| Tert-butyl ethyl ether | Ave | 0.5770 | 0.6375 | | 0.0221 | 0.0200 | 10.5 | 20.0 |
| Ethyl acetate | Ave | 0.3452 | 0.3255 | | 0.0377 | 0.0400 | -5.7 | 20.0 |
| Propionitrile | Ave | 0.0531 | 0.0421 | | 0.159 | 0.200 | -20.7* | 20.0 |
| Methacrylonitrile | Ave | 0.2706 | 0.2228 | | 0.165 | 0.200 | -17.7 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.4381 | 0.4514 | | 0.0206 | 0.0200 | 3.0 | 20.0 |
| n-Butanol | Qua | | 0.0051 | | 0.391 | 0.500 | -21.7* | 20.0 |
| Methyl methacrylate | Ave | 0.3461 | 0.3120 | | 0.0361 | 0.0400 | -9.8 | 20.0 |
| 2-Nitropropane | Qua | | 0.0620 | | 0.0365 | 0.0400 | -8.7 | 20.0 |
| 1-Chlorohexane | Ave | 0.3891 | 0.3762 | | 0.0193 | 0.0200 | -3.3 | 20.0 |
| Cyclohexanone | Qua | | 0.0078 | | 0.0914 | 0.200 | -54.3* | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 2.816 | 2.416 | | 0.0172 | 0.0200 | -14.2 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.147 | 0.8132 | | 0.0142 | 0.0200 | -29.1* | 20.0 |
| 2-Methylnaphthalene | Ave | 1.621 | 0.9593 | | 0.0237 | 0.0400 | -40.8* | 20.0 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207096/6

Matrix: Water | Lab File ID: UXX3070.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 12:17

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.28 | U | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.44 | U | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.22 | U | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.30 | U | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 0.45 | U | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.37 | U | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.44 | U | 1.0 | 0.44 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.47 | U | 5.0 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.32 | U | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.41 | U | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.82 | U | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.25 | U | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 0.23 | U | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 0.25 | U | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.19 | U | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 0.19 | U | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.27 | U | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 0.26 | U | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 0.53 | U | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.59 | U | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 0.40 | U | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 0.48 | U | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 10 | 0.99 |
| 67-64-1 | Acetone | 0.94 | U | 10 | 0.94 |
| 71-43-2 | Benzene | 0.35 | U | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 0.35 | U | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 0.56 | U | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 0.44 | U | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.38 | U | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 0.43 | U | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 0.25 | U | 1.0 | 0.25 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207096/6

Matrix: Water | Lab File ID: UXX3070.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 12:17

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 1.0 | 0.43 |
| 75-00-3 | Chloroethane | 0.32 | U | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 0.25 | U | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 0.44 | U | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.26 | U | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.46 | U | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 0.42 | U | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.50 | U | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 4.9 | U | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 0.25 | U | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 0.35 | U | 1.0 | 0.35 |
| 108-20-3 | Diisopropyl ether | 0.50 | U | 10 | 0.50 |
| 98-82-8 | Isopropylbenzene | 0.35 | U | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 0.20 | U | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 1.18 | | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.24 | U | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 0.45 | U | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 0.31 | U | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 0.40 | U | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 0.25 | U | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 0.48 | U | 1.0 | 0.48 |
| 100-42-5 | Styrene | 0.45 | U | 1.0 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.30 | U | 5.0 | 0.30 |
| 637-92-3 | Ethyl t-butyl ether | 0.23 | U | 5.0 | 0.23 |
| 127-18-4 | Tetrachloroethene | 0.31 | U | 1.0 | 0.31 |
| 108-88-3 | Toluene | 0.23 | U | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.30 | U | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.56 | U | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 0.22 | U | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 0.49 | U | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 0.41 | U | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 0.29 | U | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 0.52 | U | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 0.41 | U | 1.0 | 0.41 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207096/6

Matrix: Water | Lab File ID: UXX3070.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 12:17

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.32 | U | 1.0 | 0.32 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 82 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 82 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 75 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 82 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207343/6

Matrix: Solid | Lab File ID: 146535.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/18/2015 12:20

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.80 | U | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.78 | U | 5.0 | 0.78 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 1.1 | U | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.90 | U | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207343/6

Matrix: Solid | Lab File ID: 146535.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/18/2015 12:20

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.79 | U | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.71 | U | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.7 | U | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.74 | U | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.35 | U | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.91 | U | 5.0 | 0.91 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.80 | U | 5.0 | 0.80 |
| 108-88-3 | Toluene | 0.298 | J | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207343/6

Matrix: Solid | Lab File ID: 146535.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/18/2015 12:20

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 87 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 91 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 76 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 95 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207560/6

Matrix: Solid | Lab File ID: 146568.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/19/2015 12:29

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.80 | U | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.78 | U | 5.0 | 0.78 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 1.1 | U | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.90 | U | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207560/6

Matrix: Solid | Lab File ID: 146568.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/19/2015 12:29

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.79 | U | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.71 | U | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.7 | U | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.74 | U | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.35 | U | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.91 | U | 5.0 | 0.91 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.80 | U | 5.0 | 0.80 |
| 108-88-3 | Toluene | 0.348 | J | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207560/6

Matrix: Solid | Lab File ID: 146568.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/19/2015 12:29

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 92 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 82 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207720/8

Matrix: Solid | Lab File ID: UX80879.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/20/2015 05:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207720 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.80 | U | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.78 | U | 5.0 | 0.78 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 1.1 | U | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 6.42 | J | 20 | 5.7 |
| 71-43-2 | Benzene | 0.90 | U | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207720/8

Matrix: Solid | Lab File ID: UX80879.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/20/2015 05:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207720 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.79 | U | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.71 | U | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.7 | U | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.74 | U | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.35 | U | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.91 | U | 5.0 | 0.91 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.80 | U | 5.0 | 0.80 |
| 108-88-3 | Toluene | 0.27 | U | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207720/8

Matrix: Solid | Lab File ID: UX80879.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/20/2015 05:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207720 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 120 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 117 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 114 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 114 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207096/4

Matrix: Water | Lab File ID: UXX3068.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 11:34

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 9.46 | | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 9.11 | | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 10.0 | | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 9.72 | | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 9.27 | | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 9.08 | | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 10.0 | | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5.89 | | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 10.0 | | 1.0 | 0.44 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 6.37 | | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 9.26 | | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 8.37 | | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 9.01 | | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 9.37 | | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 9.44 | | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 9.15 | | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 9.46 | | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 9.15 | | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 7.77 | | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 19.0 | | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 9.41 | J | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 9.78 | | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 19.2 | | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 9.88 | | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 9.20 | | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 18.8 | | 10 | 0.99 |
| 67-64-1 | Acetone | 18.4 | | 10 | 0.94 |
| 71-43-2 | Benzene | 9.16 | | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 9.75 | | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 8.92 | | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 8.49 | | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 8.54 | | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 9.28 | | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 9.38 | | 1.0 | 0.25 |
| 124-48-1 | Dibromochloromethane | 9.64 | | 1.0 | 0.43 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207096/4

Matrix: Water | Lab File ID: UXX3068.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 11:34

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 8.47 | | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 9.47 | | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 8.43 | | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 9.25 | | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 9.55 | | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 9.34 | | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 9.25 | | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 6.79 | | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 9.05 | | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 73.6 | | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 9.52 | | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 6.28 | | 1.0 | 0.35 |
| 98-82-8 | Isopropylbenzene | 9.49 | | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 8.83 | | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 10.9 | | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 9.72 | | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 6.29 | | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 8.80 | | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 10.1 | | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 9.71 | | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 9.36 | | 1.0 | 0.48 |
| 100-42-5 | Styrene | 9.41 | | 1.0 | 0.45 |
| 127-18-4 | Tetrachloroethene | 9.47 | | 1.0 | 0.31 |
| 108-88-3 | Toluene | 9.21 | | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 9.18 | | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 9.95 | | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 9.58 | | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 9.63 | | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 7.50 | | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 8.72 | | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 19.4 | | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 9.70 | | 1.0 | 0.41 |
| 106-93-4 | 1,2-Dibromoethane | 9.91 | | 1.0 | 0.32 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Client Sample ID: Lab Sample ID: LCS 240-207096/4

Matrix: Water Lab File ID: UXX3068.D

Analysis Method: 8260C Date Collected:

Sample wt/vol: 5(mL) Date Analyzed: 11/17/2015 11:34

Soil Aliquot Vol: Dilution Factor: 1

Soil Extract Vol.: GC Column: DB-624 ID: 0.18(mm)

% Moisture: Level: (low/med) Low

Analysis Batch No.: 207096 Units: ug/L

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 90 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 90 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 85 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 90 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207343/5

Matrix: Solid | Lab File ID: 146534.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/18/2015 11:58

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 21.0 | | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 22.4 | | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 17.6 | | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.3 | | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 18.2 | | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 17.0 | | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 19.3 | | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 18.7 | | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 20.0 | | 5.0 | 0.78 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 18.7 | | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 18.7 | | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 18.1 | | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 18.0 | | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 19.6 | | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 16.9 | | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 17.9 | | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 17.4 | | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 17.6 | | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 22.1 | | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 26.6 | | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 12.5 | J | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 19.3 | | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 30.6 | | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 19.5 | | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 19.3 | | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 30.2 | | 20 | 1.2 |
| 67-64-1 | Acetone | 36.6 | | 20 | 5.7 |
| 71-43-2 | Benzene | 18.3 | | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 19.7 | | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 16.8 | | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 18.2 | | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 16.2 | | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 22.6 | | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 18.9 | | 5.0 | 0.53 |
| 124-48-1 | Dibromochloromethane | 18.3 | | 5.0 | 0.37 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207343/5

Matrix: Solid | Lab File ID: 146534.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/18/2015 11:58

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 17.0 | | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 19.9 | | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 13.3 | | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 20.0 | | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 16.5 | | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 19.1 | | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 20.6 | | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 16.4 | | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 18.4 | | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 135 | J | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 19.3 | | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 19.5 | | 5.0 | 0.40 |
| 98-82-8 | Isopropylbenzene | 19.6 | | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 18.0 | | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 15.3 | | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 18.8 | | 10 | 0.42 |
| 91-20-3 | Naphthalene | 16.8 | | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 18.9 | | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 19.0 | | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 20.1 | | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 18.6 | | 5.0 | 0.91 |
| 100-42-5 | Styrene | 18.2 | | 5.0 | 0.40 |
| 127-18-4 | Tetrachloroethene | 20.1 | | 5.0 | 0.80 |
| 108-88-3 | Toluene | 19.5 | | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 20.1 | | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 18.1 | | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 20.3 | | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 20.8 | | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 15.5 | | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 14.2 | | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 38.9 | | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 19.0 | | 5.0 | 0.67 |
| 106-93-4 | 1,2-Dibromoethane | 20.0 | | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-207343/5

Matrix: Solid  Lab File ID: 146534.D

Analysis Method: 8260C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/18/2015 11:58

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: DB-624 ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 207343  Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 91 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 97 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 80 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207343/7

Matrix: Solid | Lab File ID: 146536.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/18/2015 12:41

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207343 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 526-73-8 | 1,2,3-Trimethylbenzene | 18.4 | | 5.0 | 0.46 |
| 108-20-3 | Diisopropyl ether | 18.2 | | 10 | 0.59 |
| 994-05-8 | Tert-amyl methyl ether | 15.1 | | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 15.1 | | 5.0 | 0.41 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 90 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 80 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 92 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207560/5

Matrix: Solid | Lab File ID: 146567.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/19/2015 12:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 21.2 | | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 23.2 | | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 18.7 | | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 22.0 | | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 18.9 | | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 18.5 | | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 21.0 | | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 19.1 | | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 18.4 | | 5.0 | 0.78 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 19.0 | | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 19.3 | | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 18.5 | | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 18.8 | | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 20.7 | | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 17.5 | | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 18.4 | | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 18.4 | | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 18.1 | | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 23.0 | | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 30.7 | | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 13.8 | J | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 19.4 | | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 32.4 | | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 19.4 | | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 19.9 | | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 31.6 | | 20 | 1.2 |
| 67-64-1 | Acetone | 33.8 | | 20 | 5.7 |
| 71-43-2 | Benzene | 18.6 | | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 20.2 | | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 17.5 | | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 18.6 | | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 16.2 | | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 22.9 | | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 19.5 | | 5.0 | 0.53 |
| 124-48-1 | Dibromochloromethane | 18.4 | | 5.0 | 0.37 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207560/5

Matrix: Solid | Lab File ID: 146567.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/19/2015 12:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 18.5 | | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 20.8 | | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 14.3 | | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 20.5 | | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 17.9 | | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 21.3 | | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 20.9 | | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 15.4 | | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 19.3 | | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 159 | J | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 20.0 | | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 20.9 | | 5.0 | 0.40 |
| 98-82-8 | Isopropylbenzene | 20.5 | | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 18.8 | | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 15.3 | | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 19.9 | | 10 | 0.42 |
| 91-20-3 | Naphthalene | 17.7 | | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 19.4 | | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 19.3 | | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 20.3 | | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 19.4 | | 5.0 | 0.91 |
| 100-42-5 | Styrene | 18.4 | | 5.0 | 0.40 |
| 127-18-4 | Tetrachloroethene | 20.7 | | 5.0 | 0.80 |
| 108-88-3 | Toluene | 19.9 | | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 20.8 | | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 19.3 | | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 20.4 | | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 21.9 | | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 17.2 | | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 14.8 | | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 40.2 | | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 19.2 | | 5.0 | 0.67 |
| 106-93-4 | 1,2-Dibromoethane | 20.9 | | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-207560/5

Matrix: Solid  Lab File ID: 146567.D

Analysis Method: 8260C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/19/2015 12:08

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: DB-624  ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 207560  Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 92 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 81 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 98 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207560/7

Matrix: Solid | Lab File ID: 146569.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/19/2015 12:51

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 526-73-8 | 1,2,3-Trimethylbenzene | 18.5 | | 5.0 | 0.46 |
| 108-20-3 | Diisopropyl ether | 18.0 | | 10 | 0.59 |
| 994-05-8 | Tert-amyl methyl ether | 14.9 | | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 15.7 | | 5.0 | 0.41 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 95 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 83 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207720/6

Matrix: Solid | Lab File ID: UX80877.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/20/2015 04:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207720 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 49.7 | | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 49.6 | | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 49.8 | | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 51.8 | | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 48.9 | | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 48.3 | | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 52.2 | | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 45.0 | | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 52.8 | | 5.0 | 0.78 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 46.3 | | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 49.4 | | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 48.6 | | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 46.6 | | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 49.9 | | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 47.1 | | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 45.9 | | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 46.0 | | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 45.5 | | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 44.2 | | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 98.6 | | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | NC | | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 48.0 | | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 102 | | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 45.9 | | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 50.3 | | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 107 | | 20 | 1.2 |
| 67-64-1 | Acetone | 111 | | 20 | 5.7 |
| 71-43-2 | Benzene | 46.8 | | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 46.6 | | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 49.8 | | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 56.1 | | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 50.4 | | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 53.3 | | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 45.5 | | 5.0 | 0.53 |
| 124-48-1 | Dibromochloromethane | 52.7 | | 5.0 | 0.37 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207720/6

Matrix: Solid | Lab File ID: UX80877.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/20/2015 04:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207720 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 49.4 | | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 49.5 | | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 50.8 | | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 48.3 | | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 48.3 | | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 49.9 | | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 49.5 | | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 52.4 | | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 50.4 | | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 440 | | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 45.5 | | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 48.6 | | 5.0 | 0.40 |
| 98-82-8 | Isopropylbenzene | 50.1 | | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 47.8 | | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 52.3 | | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 46.2 | | 10 | 0.42 |
| 91-20-3 | Naphthalene | 52.0 | | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 50.0 | | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 48.6 | | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 47.8 | | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 51.0 | | 5.0 | 0.91 |
| 100-42-5 | Styrene | 45.9 | | 5.0 | 0.40 |
| 127-18-4 | Tetrachloroethene | 45.4 | | 5.0 | 0.80 |
| 108-88-3 | Toluene | 45.5 | | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 48.7 | | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 42.9 | | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 48.9 | | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 50.8 | | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 39.2 | | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 49.9 | | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 94.0 | | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 55.3 | | 5.0 | 0.67 |
| 106-93-4 | 1,2-Dibromoethane | 48.6 | | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-207720/6

Matrix: Solid  Lab File ID: UX80877.D

Analysis Method: 8260C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/20/2015 04:26

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: DB-624 ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 207720  Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 104 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 105 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 96 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 100 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 200-30727-B-1 MS

Matrix: Water | Lab File ID: UXX3079.D

Analysis Method: 8260C | Date Collected: 11/11/2015 14:30

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 15:35

Soil Aliquot Vol: | Dilution Factor: 50

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 486 | | 50 | 14 |
| 71-55-6 | 1,1,1-Trichloroethane | 466 | | 50 | 22 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 484 | | 50 | 11 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 502 | | 50 | 23 |
| 75-34-3 | 1,1-Dichloroethane | 477 | | 50 | 15 |
| 75-35-4 | 1,1-Dichloroethene | 485 | | 50 | 23 |
| 563-58-6 | 1,1-Dichloropropene | 526 | | 50 | 21 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 409 | | 50 | 19 |
| 96-18-4 | 1,2,3-Trichloropropane | 498 | | 50 | 22 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 426 | | 50 | 16 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 477 | | 50 | 21 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 381 | | 100 | 41 |
| 95-50-1 | 1,2-Dichlorobenzene | 475 | | 50 | 13 |
| 107-06-2 | 1,2-Dichloroethane | 489 | | 50 | 12 |
| 78-87-5 | 1,2-Dichloropropane | 483 | | 50 | 13 |
| 541-73-1 | 1,3-Dichlorobenzene | 468 | | 50 | 9.5 |
| 142-28-9 | 1,3-Dichloropropane | 483 | | 50 | 9.5 |
| 106-46-7 | 1,4-Dichlorobenzene | 472 | | 50 | 14 |
| 594-20-7 | 2,2-Dichloropropane | 402 | | 50 | 13 |
| 78-93-3 | 2-Butanone (MEK) | 953 | | 500 | 27 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 278 | J | 500 | 30 |
| 95-49-8 | 2-Chlorotoluene | 484 | | 50 | 20 |
| 591-78-6 | 2-Hexanone | 954 | | 500 | 24 |
| 106-43-4 | 4-Chlorotoluene | 501 | | 50 | 15 |
| 99-87-6 | p-Isopropyltoluene | 475 | | 50 | 22 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 957 | | 500 | 50 |
| 67-64-1 | Acetone | 998 | | 500 | 47 |
| 71-43-2 | Benzene | 471 | | 50 | 18 |
| 108-86-1 | Bromobenzene | 486 | | 50 | 18 |
| 75-25-2 | Bromoform | 419 | | 50 | 28 |
| 74-83-9 | Bromomethane | 402 | | 50 | 22 |
| 75-15-0 | Carbon disulfide | 465 | | 50 | 19 |
| 56-23-5 | Carbon tetrachloride | 475 | | 50 | 22 |
| 108-90-7 | Chlorobenzene | 482 | | 50 | 13 |
| 124-48-1 | Dibromochloromethane | 458 | | 50 | 22 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 200-30727-B-1 MS

Matrix: Water | Lab File ID: UXX3079.D

Analysis Method: 8260C | Date Collected: 11/11/2015 14:30

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 15:35

Soil Aliquot Vol: | Dilution Factor: 50

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 426 | | 50 | 16 |
| 67-66-3 | Chloroform | 488 | | 50 | 13 |
| 74-87-3 | Chloromethane | 356 | | 50 | 22 |
| 156-59-2 | cis-1,2-Dichloroethene | 1660 | | 50 | 13 |
| 10061-01-5 | cis-1,3-Dichloropropene | 470 | | 50 | 23 |
| 74-95-3 | Dibromomethane | 479 | | 50 | 21 |
| 74-97-5 | Bromochloromethane | 481 | | 50 | 25 |
| 75-71-8 | Dichlorodifluoromethane | 360 | | 50 | 16 |
| 75-27-4 | Bromodichloromethane | 458 | | 50 | 15 |
| 100-41-4 | Ethylbenzene | 489 | | 50 | 13 |
| 87-68-3 | Hexachlorobutadiene | 395 | | 50 | 18 |
| 98-82-8 | Isopropylbenzene | 498 | | 50 | 18 |
| 1634-04-4 | Methyl tert-butyl ether | 443 | | 50 | 10 |
| 75-09-2 | Methylene Chloride | 578 | | 50 | 17 |
| 179601-23-1 | m-Xylene & p-Xylene | 509 | | 100 | 12 |
| 91-20-3 | Naphthalene | 498 | | 50 | 23 |
| 104-51-8 | n-Butylbenzene | 473 | | 50 | 16 |
| 103-65-1 | N-Propylbenzene | 516 | | 50 | 20 |
| 95-47-6 | o-Xylene | 501 | | 50 | 13 |
| 135-98-8 | sec-Butylbenzene | 489 | | 50 | 24 |
| 100-42-5 | Styrene | 487 | | 50 | 23 |
| 127-18-4 | Tetrachloroethene | 1140 | | 50 | 16 |
| 108-88-3 | Toluene | 483 | | 50 | 12 |
| 156-60-5 | trans-1,2-Dichloroethene | 493 | | 50 | 15 |
| 10061-02-6 | trans-1,3-Dichloropropene | 474 | | 50 | 28 |
| 79-01-6 | Trichloroethene | 1110 | | 50 | 11 |
| 75-69-4 | Trichlorofluoromethane | 455 | | 50 | 25 |
| 108-05-4 | Vinyl acetate | 376 | | 100 | 21 |
| 75-01-4 | Vinyl chloride | 693 | | 50 | 15 |
| 1330-20-7 | Xylenes, Total | 1010 | | 100 | 26 |
| 98-06-6 | tert-Butylbenzene | 484 | | 50 | 21 |
| 106-93-4 | 1,2-Dibromoethane | 494 | | 50 | 16 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: 200-30727-B-1 MS

Matrix: Water  Lab File ID: UXX3079.D

Analysis Method: 8260C  Date Collected: 11/11/2015 14:30

Sample wt/vol: 5(mL)  Date Analyzed: 11/17/2015 15:35

Soil Aliquot Vol:  Dilution Factor: 50

Soil Extract Vol.:  GC Column: DB-624 ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 207096  Units: ug/L

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 84 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 81 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 80 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 81 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57921-A-4-J MS

Matrix: Solid | Lab File ID: 146590.D

Analysis Method: 8260C | Date Collected: 11/10/2015 00:00

Sample wt/vol: 5.16(g) | Date Analyzed: 11/19/2015 20:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 4.3 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 71-55-6 | 1,1,1-Trichloroethane | 20.0 | | 5.1 | 0.66 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.1 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 21.1 | | 5.1 | 0.43 |
| 75-34-3 | 1,1-Dichloroethane | 16.3 | | 5.1 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 24.7 | | 5.1 | 0.81 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 9.56 | | 5.1 | 0.38 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 6.61 | J | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 11.6 | | 5.1 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 18.3 | | 5.1 | 0.47 |
| 78-87-5 | 1,2-Dichloropropane | 15.3 | | 5.1 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 10.5 | | 5.1 | 0.70 |
| 106-46-7 | 1,4-Dichlorobenzene | 10.3 | | 5.1 | 0.28 |
| 78-93-3 | 2-Butanone (MEK) | 32.2 | | 20 | 1.1 |
| 591-78-6 | 2-Hexanone | 29.9 | | 20 | 0.69 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 28.0 | | 20 | 1.2 |
| 67-64-1 | Acetone | 78.3 | | 20 | 5.8 |
| 71-43-2 | Benzene | 16.0 | | 5.1 | 0.91 |
| 75-25-2 | Bromoform | 14.0 | | 5.1 | 0.24 |
| 74-83-9 | Bromomethane | 12.4 | | 5.1 | 0.43 |
| 75-15-0 | Carbon disulfide | 15.1 | | 5.1 | 0.60 |
| 56-23-5 | Carbon tetrachloride | 19.5 | | 5.1 | 0.66 |
| 108-90-7 | Chlorobenzene | 14.1 | | 5.1 | 0.54 |
| 124-48-1 | Dibromochloromethane | 14.4 | | 5.1 | 0.37 |
| 75-00-3 | Chloroethane | 15.8 | | 5.1 | 0.45 |
| 67-66-3 | Chloroform | 18.3 | | 5.1 | 0.36 |
| 74-87-3 | Chloromethane | 12.2 | | 5.1 | 0.80 |
| 156-59-2 | cis-1,2-Dichloroethene | 17.3 | | 5.1 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 9.86 | | 5.1 | 0.72 |
| 75-71-8 | Dichlorodifluoromethane | 13.8 | | 5.1 | 0.28 |
| 75-27-4 | Bromodichloromethane | 14.2 | | 5.1 | 0.29 |
| 100-41-4 | Ethylbenzene | 14.0 | | 5.1 | 0.32 |
| 98-82-8 | Isopropylbenzene | 13.3 | | 5.1 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 17.4 | | 5.1 | 0.44 |
| 75-09-2 | Methylene Chloride | 16.2 | | 5.1 | 0.75 |
| 179601-23-1 | m-Xylene & p-Xylene | 13.6 | | 10 | 0.43 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57921-A-4-J MS

Matrix: Solid | Lab File ID: 146590.D

Analysis Method: 8260C | Date Collected: 11/10/2015 00:00

Sample wt/vol: 5.16(g) | Date Analyzed: 11/19/2015 20:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 4.3 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 95-47-6 | o-Xylene | 14.4 | | 5.1 | 0.58 |
| 100-42-5 | Styrene | 12.7 | | 5.1 | 0.40 |
| 127-18-4 | Tetrachloroethene | 15.3 | | 5.1 | 0.81 |
| 108-88-3 | Toluene | 15.9 | | 5.1 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 17.3 | | 5.1 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 12.3 | | 5.1 | 0.29 |
| 79-01-6 | Trichloroethene | 31.6 | | 5.1 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 18.7 | | 5.1 | 0.29 |
| 75-01-4 | Vinyl chloride | 13.6 | | 5.1 | 0.30 |
| 1330-20-7 | Xylenes, Total | 28.0 | | 10 | 0.55 |
| 106-93-4 | 1,2-Dibromoethane | 17.1 | | 5.1 | 0.54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 90 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 52 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 83 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57972-C-3-C MS

Matrix: Solid | Lab File ID: UX80889.D

Analysis Method: 8260C | Date Collected: 11/16/2015 15:50

Sample wt/vol: 5.28(g) | Date Analyzed: 11/20/2015 08:41

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 12.6 | Level: (low/med) Low

Analysis Batch No.: 207720 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 67-64-1 | Acetone | 37.4 | | 22 | 6.2 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 101 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 98 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 200-30727-B-1 MSD

Matrix: Water | Lab File ID: UXX3080.D

Analysis Method: 8260C | Date Collected: 11/11/2015 14:30

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 15:57

Soil Aliquot Vol: | Dilution Factor: 50

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 467 | | 50 | 14 |
| 71-55-6 | 1,1,1-Trichloroethane | 479 | | 50 | 22 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 485 | | 50 | 11 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 518 | | 50 | 23 |
| 75-34-3 | 1,1-Dichloroethane | 479 | | 50 | 15 |
| 75-35-4 | 1,1-Dichloroethene | 504 | | 50 | 23 |
| 563-58-6 | 1,1-Dichloropropene | 514 | | 50 | 21 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 414 | | 50 | 19 |
| 96-18-4 | 1,2,3-Trichloropropane | 491 | | 50 | 22 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 425 | | 50 | 16 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 470 | | 50 | 21 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 387 | | 100 | 41 |
| 95-50-1 | 1,2-Dichlorobenzene | 465 | | 50 | 13 |
| 107-06-2 | 1,2-Dichloroethane | 487 | | 50 | 12 |
| 78-87-5 | 1,2-Dichloropropane | 486 | | 50 | 13 |
| 541-73-1 | 1,3-Dichlorobenzene | 469 | | 50 | 9.5 |
| 142-28-9 | 1,3-Dichloropropane | 483 | | 50 | 9.5 |
| 106-46-7 | 1,4-Dichlorobenzene | 465 | | 50 | 14 |
| 594-20-7 | 2,2-Dichloropropane | 410 | | 50 | 13 |
| 78-93-3 | 2-Butanone (MEK) | 997 | | 500 | 27 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 254 | J | 500 | 30 |
| 95-49-8 | 2-Chlorotoluene | 485 | | 50 | 20 |
| 591-78-6 | 2-Hexanone | 985 | | 500 | 24 |
| 106-43-4 | 4-Chlorotoluene | 486 | | 50 | 15 |
| 99-87-6 | p-Isopropyltoluene | 477 | | 50 | 22 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 965 | | 500 | 50 |
| 67-64-1 | Acetone | 1070 | | 500 | 47 |
| 71-43-2 | Benzene | 469 | | 50 | 18 |
| 108-86-1 | Bromobenzene | 494 | | 50 | 18 |
| 75-25-2 | Bromoform | 416 | | 50 | 28 |
| 74-83-9 | Bromomethane | 420 | | 50 | 22 |
| 75-15-0 | Carbon disulfide | 455 | | 50 | 19 |
| 56-23-5 | Carbon tetrachloride | 485 | | 50 | 22 |
| 108-90-7 | Chlorobenzene | 477 | | 50 | 13 |
| 124-48-1 | Dibromochloromethane | 464 | | 50 | 22 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 200-30727-B-1 MSD

Matrix: Water | Lab File ID: UXX3080.D

Analysis Method: 8260C | Date Collected: 11/11/2015 14:30

Sample wt/vol: 5(mL) | Date Analyzed: 11/17/2015 15:57

Soil Aliquot Vol: | Dilution Factor: 50

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 207096 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 432 | | 50 | 16 |
| 67-66-3 | Chloroform | 492 | | 50 | 13 |
| 74-87-3 | Chloromethane | 379 | | 50 | 22 |
| 156-59-2 | cis-1,2-Dichloroethene | 1660 | | 50 | 13 |
| 10061-01-5 | cis-1,3-Dichloropropene | 483 | | 50 | 23 |
| 74-95-3 | Dibromomethane | 478 | | 50 | 21 |
| 74-97-5 | Bromochloromethane | 467 | | 50 | 25 |
| 75-71-8 | Dichlorodifluoromethane | 354 | | 50 | 16 |
| 75-27-4 | Bromodichloromethane | 462 | | 50 | 15 |
| 100-41-4 | Ethylbenzene | 489 | | 50 | 13 |
| 87-68-3 | Hexachlorobutadiene | 410 | | 50 | 18 |
| 98-82-8 | Isopropylbenzene | 479 | | 50 | 18 |
| 1634-04-4 | Methyl tert-butyl ether | 457 | | 50 | 10 |
| 75-09-2 | Methylene Chloride | 579 | | 50 | 17 |
| 179601-23-1 | m-Xylene & p-Xylene | 488 | | 100 | 12 |
| 91-20-3 | Naphthalene | 431 | | 50 | 23 |
| 104-51-8 | n-Butylbenzene | 463 | | 50 | 16 |
| 103-65-1 | N-Propylbenzene | 510 | | 50 | 20 |
| 95-47-6 | o-Xylene | 492 | | 50 | 13 |
| 135-98-8 | sec-Butylbenzene | 480 | | 50 | 24 |
| 100-42-5 | Styrene | 464 | | 50 | 23 |
| 127-18-4 | Tetrachloroethene | 1160 | | 50 | 16 |
| 108-88-3 | Toluene | 478 | | 50 | 12 |
| 156-60-5 | trans-1,2-Dichloroethene | 486 | | 50 | 15 |
| 10061-02-6 | trans-1,3-Dichloropropene | 494 | | 50 | 28 |
| 79-01-6 | Trichloroethene | 1110 | | 50 | 11 |
| 75-69-4 | Trichlorofluoromethane | 479 | | 50 | 25 |
| 108-05-4 | Vinyl acetate | 383 | | 100 | 21 |
| 75-01-4 | Vinyl chloride | 716 | | 50 | 15 |
| 1330-20-7 | Xylenes, Total | 980 | | 100 | 26 |
| 98-06-6 | tert-Butylbenzene | 485 | | 50 | 21 |
| 106-93-4 | 1,2-Dibromoethane | 485 | | 50 | 16 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: 200-30727-B-1 MSD

Matrix: Water  Lab File ID: UXX3080.D

Analysis Method: 8260C  Date Collected: 11/11/2015 14:30

Sample wt/vol: 5(mL)  Date Analyzed: 11/17/2015 15:57

Soil Aliquot Vol:  Dilution Factor: 50

Soil Extract Vol.:  GC Column: DB-624 ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 207096  Units: ug/L

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 84 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 84 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 80 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 80 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57921-A-4-K MSD

Matrix: Solid | Lab File ID: 146591.D

Analysis Method: 8260C | Date Collected: 11/10/2015 00:00

Sample wt/vol: 5.24(g) | Date Analyzed: 11/19/2015 20:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 4.3 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 71-55-6 | 1,1,1-Trichloroethane | 20.2 | | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 17.4 | | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 16.3 | | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 22.7 | | 5.0 | 0.80 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 10.9 | | 5.0 | 0.38 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 5.77 | J | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 13.2 | | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 17.9 | | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 15.8 | | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 12.7 | | 5.0 | 0.69 |
| 106-46-7 | 1,4-Dichlorobenzene | 11.8 | | 5.0 | 0.28 |
| 78-93-3 | 2-Butanone (MEK) | 28.6 | | 20 | 1.1 |
| 591-78-6 | 2-Hexanone | 28.2 | | 20 | 0.68 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 27.1 | | 20 | 1.2 |
| 67-64-1 | Acetone | 71.4 | | 20 | 5.7 |
| 71-43-2 | Benzene | 16.5 | | 5.0 | 0.90 |
| 75-25-2 | Bromoform | 13.5 | | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 11.7 | | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 15.5 | | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 20.0 | | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 14.9 | | 5.0 | 0.53 |
| 124-48-1 | Dibromochloromethane | 13.5 | | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 16.7 | | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 18.6 | | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 12.1 | | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 17.6 | | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 9.38 | | 5.0 | 0.71 |
| 75-71-8 | Dichlorodifluoromethane | 13.7 | | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 13.6 | | 5.0 | 0.29 |
| 100-41-4 | Ethylbenzene | 15.4 | | 5.0 | 0.32 |
| 98-82-8 | Isopropylbenzene | 15.4 | | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 17.2 | | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 17.4 | | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 14.9 | | 10 | 0.42 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57921-A-4-K MSD

Matrix: Solid | Lab File ID: 146591.D

Analysis Method: 8260C | Date Collected: 11/10/2015 00:00

Sample wt/vol: 5.24(g) | Date Analyzed: 11/19/2015 20:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 4.3 | Level: (low/med) Low

Analysis Batch No.: 207560 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 95-47-6 | o-Xylene | 15.7 | | 5.0 | 0.57 |
| 100-42-5 | Styrene | 13.4 | | 5.0 | 0.40 |
| 127-18-4 | Tetrachloroethene | 16.8 | | 5.0 | 0.80 |
| 108-88-3 | Toluene | 16.7 | | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 17.6 | | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 11.4 | | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 32.5 | | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 19.5 | | 5.0 | 0.29 |
| 75-01-4 | Vinyl chloride | 13.7 | | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 30.6 | | 10 | 0.54 |
| 106-93-4 | 1,2-Dibromoethane | 16.6 | | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 90 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 49 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 84 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57972-C-3-D MSD

Matrix: Solid | Lab File ID: UX80890.D

Analysis Method: 8260C | Date Collected: 11/16/2015 15:50

Sample wt/vol: 5.04(g) | Date Analyzed: 11/20/2015 09:02

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 12.6 | Level: (low/med) Low

Analysis Batch No.: 207720 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 67-64-1 | Acetone | 46.4 | | 23 | 6.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 101 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 100 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 95 | | 58-123 |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX10 Start Date: 11/16/2015 09:00

Analysis Batch Number: 206893 End Date: 11/16/2015 19:02

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-206893/1 | | 11/16/2015 09:00 | 1 | BFB7317.D | DB-624 0.18(mm) |
| STD8260 240-206893/2 IC | | 11/16/2015 09:23 | 1 | UXX3039.D | DB-624 0.18(mm) |
| STD8260 240-206893/3 IC | | 11/16/2015 09:45 | 1 | UXX3040.D | DB-624 0.18(mm) |
| STD8260 240-206893/4 ICIS | | 11/16/2015 10:07 | 1 | UXX3041.D | DB-624 0.18(mm) |
| STD8260 240-206893/5 IC | | 11/16/2015 10:29 | 1 | UXX3042.D | DB-624 0.18(mm) |
| STD8260 240-206893/6 IC | | 11/16/2015 10:51 | 1 | UXX3043.D | DB-624 0.18(mm) |
| STD8260 240-206893/7 IC | | 11/16/2015 11:13 | 1 | UXX3044.D | DB-624 0.18(mm) |
| STD8260 240-206893/8 IC | | 11/16/2015 11:35 | 1 | UXX3045.D | DB-624 0.18(mm) |
| ICV 240-206893/9 | | 11/16/2015 11:57 | 1 | | DB-624 0.18(mm) |
| STDA9 240-206893/10 IC | | 11/16/2015 12:19 | 1 | UXX3047.D | DB-624 0.18(mm) |
| STDA9 240-206893/11 IC | | 11/16/2015 12:43 | 1 | UXX3048.D | DB-624 0.18(mm) |
| STDA9 240-206893/12 IC | | 11/16/2015 13:05 | 1 | UXX3049.D | DB-624 0.18(mm) |
| STDA9 240-206893/13 IC | | 11/16/2015 13:27 | 1 | UXX3050.D | DB-624 0.18(mm) |
| STDA9 240-206893/14 IC | | 11/16/2015 13:49 | 1 | UXX3051.D | DB-624 0.18(mm) |
| STDA9 240-206893/15 IC | | 11/16/2015 14:10 | 1 | UXX3052.D | DB-624 0.18(mm) |
| ICV 240-206893/16 | | 11/16/2015 14:33 | 1 | UXX3053.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/16/2015 15:43 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/16/2015 16:27 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/16/2015 16:49 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/16/2015 17:11 | 1 | | DB-624 0.18(mm) |
| 240-57093-A-1 MDLV | | 11/16/2015 17:33 | 1 | | DB-624 0.18(mm) |
| 240-57093-A-1 MDLV | | 11/16/2015 17:56 | 1 | | DB-624 0.18(mm) |
| 240-57093-A-1 MDLV | | 11/16/2015 18:18 | 1 | | DB-624 0.18(mm) |
| 240-57093-A-1 MDLV | | 11/16/2015 18:39 | 1 | | DB-624 0.18(mm) |
| 240-57093-A-1 MDLV | | 11/16/2015 19:02 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX10 Start Date: 11/17/2015 10:26

Analysis Batch Number: 207096 End Date: 11/17/2015 15:57

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-207096/1 | | 11/17/2015 10:26 | 1 | BFB7318.D | DB-624 0.18(mm) |
| CCVIS 240-207096/2 | | 11/17/2015 10:50 | 1 | UXX3066.D | DB-624 0.18(mm) |
| CCV 240-207096/3 | | 11/17/2015 11:11 | 1 | UXX3067.D | DB-624 0.18(mm) |
| LCS 240-207096/4 | | 11/17/2015 11:34 | 1 | UXX3068.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/17/2015 11:56 | 1 | | DB-624 0.18(mm) |
| MB 240-207096/6 | | 11/17/2015 12:17 | 1 | UXX3070.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/17/2015 12:39 | 50 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/17/2015 13:23 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/17/2015 13:45 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/17/2015 14:07 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/17/2015 14:29 | 1 | | DB-624 0.18(mm) |
| 240-57930-12 | TB-111315 | 11/17/2015 14:51 | 1 | UXX3077.D | DB-624 0.18(mm) |
| 200-30727-B-1 MS | | 11/17/2015 15:35 | 50 | UXX3079.D | DB-624 0.18(mm) |
| 200-30727-B-1 MSD | | 11/17/2015 15:57 | 50 | UXX3080.D | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX14 Start Date: 08/13/2015 13:54

Analysis Batch Number: 193150 End Date: 08/13/2015 23:12

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-193150/28 | | 08/13/2015 13:54 | 1 | BFB14045.D | DB-624 0.18(mm) |
| IC 240-193150/3 | | 08/13/2015 14:36 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/4 | | 08/13/2015 14:58 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/5 | | 08/13/2015 15:20 | 1 | | DB-624 0.18(mm) |
| ICIS 240-193150/6 | | 08/13/2015 15:41 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/7 | | 08/13/2015 16:03 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/8 | | 08/13/2015 16:24 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/9 | | 08/13/2015 16:45 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/10 | | 08/13/2015 17:07 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/11 | | 08/13/2015 17:28 | 1 | | DB-624 0.18(mm) |
| IC 240-193150/12 | | 08/13/2015 17:50 | 1 | 144576.D | DB-624 0.18(mm) |
| IC 240-193150/13 | | 08/13/2015 18:11 | 1 | 144577.D | DB-624 0.18(mm) |
| IC 240-193150/14 | | 08/13/2015 18:33 | 1 | 144578.D | DB-624 0.18(mm) |
| IC 240-193150/15 | | 08/13/2015 18:54 | 1 | 144579.D | DB-624 0.18(mm) |
| IC 240-193150/16 | | 08/13/2015 19:16 | 1 | 144580.D | DB-624 0.18(mm) |
| IC 240-193150/17 | | 08/13/2015 19:37 | 1 | 144581.D | DB-624 0.18(mm) |
| IC 240-193150/18 | | 08/13/2015 19:59 | 1 | 144582.D | DB-624 0.18(mm) |
| IC 240-193150/19 | | 08/13/2015 20:20 | 1 | 144583.D | DB-624 0.18(mm) |
| IC 240-193150/20 | | 08/13/2015 20:42 | 1 | 144584.D | DB-624 0.18(mm) |
| ICV 240-193150/21 | | 08/13/2015 21:03 | 1 | | DB-624 0.18(mm) |
| ICV 240-193150/22 | | 08/13/2015 21:25 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 21:46 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 22:08 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 22:29 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 22:50 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 08/13/2015 23:12 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX14 Start Date: 09/02/2015 13:49

Analysis Batch Number: 196026 End Date: 09/02/2015 19:33

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-196026/1 | | 09/02/2015 13:49 | 1 | BFB14070.D | DB-624 0.18(mm) |
| IC 240-196026/3 | | 09/02/2015 14:32 | 1 | 144935.D | DB-624 0.18(mm) |
| IC 240-196026/4 | | 09/02/2015 14:54 | 1 | 144936.D | DB-624 0.18(mm) |
| IC 240-196026/5 | | 09/02/2015 15:15 | 1 | 144937.D | DB-624 0.18(mm) |
| ICIS 240-196026/6 | | 09/02/2015 15:37 | 1 | 144938.D | DB-624 0.18(mm) |
| IC 240-196026/7 | | 09/02/2015 15:58 | 1 | 144939.D | DB-624 0.18(mm) |
| IC 240-196026/8 | | 09/02/2015 16:20 | 1 | 144940.D | DB-624 0.18(mm) |
| IC 240-196026/9 | | 09/02/2015 16:41 | 1 | 144941.D | DB-624 0.18(mm) |
| IC 240-196026/10 | | 09/02/2015 17:03 | 1 | 144942.D | DB-624 0.18(mm) |
| IC 240-196026/11 | | 09/02/2015 17:24 | 1 | 144943.D | DB-624 0.18(mm) |
| CCV 240-196026/12 | | 09/02/2015 17:46 | 1 | | DB-624 0.18(mm) |
| ICV 240-196026/13 | | 09/02/2015 18:07 | 1 | 144945.D | DB-624 0.18(mm) |
| ZZZZZ | | 09/02/2015 18:29 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 09/02/2015 18:50 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 09/02/2015 19:12 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 09/02/2015 19:33 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX14 Start Date: 11/18/2015 10:30

Analysis Batch Number: 207343 End Date: 11/18/2015 21:59

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-207343/1 | | 11/18/2015 10:30 | 1 | BFB14134.D | DB-624 0.18(mm) |
| CCVIS 240-207343/3 | | 11/18/2015 11:15 | 1 | 146532.D | DB-624 0.18(mm) |
| CCV 240-207343/4 | | 11/18/2015 11:37 | 1 | 146533.D | DB-624 0.18(mm) |
| LCS 240-207343/5 | | 11/18/2015 11:58 | 1 | 146534.D | DB-624 0.18(mm) |
| MB 240-207343/6 | | 11/18/2015 12:20 | 1 | 146535.D | DB-624 0.18(mm) |
| LCS 240-207343/7 | | 11/18/2015 12:41 | 1 | 146536.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 13:03 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 13:24 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 13:45 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 14:07 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 15:33 | 1 | | DB-624 0.18(mm) |
| 240-57930-1 | F-VS-UST-P-1 | 11/18/2015 18:03 | 1 | 146551.D | DB-624 0.18(mm) |
| 240-57930-3 | F-VS-UST-P-3 | 11/18/2015 18:25 | 1 | 146552.D | DB-624 0.18(mm) |
| 240-57930-5 | F-VB-UST-P-1 | 11/18/2015 19:51 | 1 | 146556.D | DB-624 0.18(mm) |
| 240-57930-6 | F-VB-UST-P-2 | 11/18/2015 20:12 | 1 | 146557.D | DB-624 0.18(mm) |
| 240-57930-7 | F-VB-UST-P-3 | 11/18/2015 20:33 | 1 | 146558.D | DB-624 0.18(mm) |
| 240-57930-8 | F-VB-UST-P-4 | 11/18/2015 20:55 | 1 | 146559.D | DB-624 0.18(mm) |
| 240-57930-9 | F-VB-UST-P-5 | 11/18/2015 21:16 | 1 | 146560.D | DB-624 0.18(mm) |
| 240-57930-10 | F-VB-UST-P-6 | 11/18/2015 21:38 | 1 | 146561.D | DB-624 0.18(mm) |
| 240-57930-11 | F-VS-UST-P-DUP | 11/18/2015 21:59 | 1 | 146562.D | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX14 Start Date: 11/19/2015 10:42

Analysis Batch Number: 207560 End Date: 11/19/2015 22:09

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-207560/11 | | 11/19/2015 10:42 | 1 | BFB14136.D | DB-624 0.18(mm) |
| CCVIS 240-207560/3 | | 11/19/2015 11:25 | 1 | 146565.D | DB-624 0.18(mm) |
| CCV 240-207560/4 | | 11/19/2015 11:46 | 1 | 146566.D | DB-624 0.18(mm) |
| LCS 240-207560/5 | | 11/19/2015 12:08 | 1 | 146567.D | DB-624 0.18(mm) |
| MB 240-207560/6 | | 11/19/2015 12:29 | 1 | 146568.D | DB-624 0.18(mm) |
| LCS 240-207560/7 | | 11/19/2015 12:51 | 1 | 146569.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/19/2015 13:12 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/19/2015 13:34 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/19/2015 16:47 | 1 | | DB-624 0.18(mm) |
| 240-57930-4 | F-VS-UST-P-4 | 11/19/2015 17:08 | 1 | 146581.D | DB-624 0.18(mm) |
| 240-57930-2 | F-VS-UST-P-2 | 11/19/2015 17:52 | 1 | 146583.D | DB-624 0.18(mm) |
| 240-57930-2 | F-VS-UST-P-2 | 11/19/2015 18:13 | 1 | 146584.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/19/2015 20:00 | 1 | | DB-624 0.18(mm) |
| 240-57921-A-4-J MS | | 11/19/2015 20:22 | 1 | 146590.D | DB-624 0.18(mm) |
| 240-57921-A-4-K MSD | | 11/19/2015 20:43 | 1 | 146591.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/19/2015 21:48 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/19/2015 22:09 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX8 Start Date: 07/15/2015 00:55

Analysis Batch Number: 189147 End Date: 07/15/2015 10:22

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-189147/1 | | 07/15/2015 00:55 | 1 | BFB8851.D | DB-624 0.18(mm) |
| CCVIS 240-189147/3 | | 07/15/2015 01:30 | 1 | | DB-624 0.18(mm) |
| STDA9 240-189147/4 IC | | 07/15/2015 01:51 | 1 | UX88545.D | DB-624 0.18(mm) |
| STDA9 240-189147/5 IC | | 07/15/2015 02:12 | 1 | UX88546.D | DB-624 0.18(mm) |
| STDA9 240-189147/6 IC | | 07/15/2015 02:34 | 1 | UX88547.D | DB-624 0.18(mm) |
| STDA9 240-189147/7 IC | | 07/15/2015 02:55 | 1 | UX88548.D | DB-624 0.18(mm) |
| STDA9 240-189147/8 IC | | 07/15/2015 03:16 | 1 | UX88549.D | DB-624 0.18(mm) |
| STDA9 240-189147/9 IC | | 07/15/2015 03:38 | 1 | UX88550.D | DB-624 0.18(mm) |
| ICV 240-189147/10 | | 07/15/2015 03:59 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 07/15/2015 04:20 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 07/15/2015 05:03 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 07/15/2015 09:40 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 07/15/2015 10:01 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 07/15/2015 10:22 | 12.5 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX8 Start Date: 07/30/2015 02:03

Analysis Batch Number: 191232 End Date: 07/30/2015 06:29

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-191232/1 | | 07/30/2015 02:03 | 1 | BFB8863.D | DB-624 0.18(mm) |
| STD8260 240-191232/3 IC | | 07/30/2015 02:35 | 1 | UX88718.D | DB-624 0.18(mm) |
| STD8260 240-191232/4 IC | | 07/30/2015 02:56 | 1 | UX88719.D | DB-624 0.18(mm) |
| STD8260 240-191232/5 IC | | 07/30/2015 03:17 | 1 | UX88720.D | DB-624 0.18(mm) |
| ICIS 240-191232/6 | | 07/30/2015 03:39 | 1 | UX88721.D | DB-624 0.18(mm) |
| STD8260 240-191232/7 IC | | 07/30/2015 04:00 | 1 | UX88722.D | DB-624 0.18(mm) |
| STD8260 240-191232/8 IC | | 07/30/2015 04:21 | 1 | UX88723.D | DB-624 0.18(mm) |
| STD8260 240-191232/9 IC | | 07/30/2015 04:42 | 1 | UX88724.D | DB-624 0.18(mm) |
| STD8260 240-191232/10 IC | | 07/30/2015 05:04 | 1 | UX88725.D | DB-624 0.18(mm) |
| ICV 240-191232/13 | | 07/30/2015 06:07 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 07/30/2015 06:29 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX8 Start Date: 11/17/2015 03:23

Analysis Batch Number: 206828 End Date: 11/17/2015 07:41

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-206828/1 | | 11/17/2015 03:23 | 1 | BFB8945.D | DB-624 0.18(mm) |
| STD8260 240-206828/3 IC | | 11/17/2015 03:57 | 1 | UX80794.D | DB-624 0.18(mm) |
| STD8260 240-206828/4 IC | | 11/17/2015 04:18 | 1 | UX80795.D | DB-624 0.18(mm) |
| STD8260 240-206828/5 IC | | 11/17/2015 04:39 | 1 | UX80796.D | DB-624 0.18(mm) |
| ICIS 240-206828/6 | | 11/17/2015 05:01 | 1 | UX80797.D | DB-624 0.18(mm) |
| STD8260 240-206828/7 IC | | 11/17/2015 05:34 | 1 | UX80798.D | DB-624 0.18(mm) |
| STD8260 240-206828/8 IC | | 11/17/2015 05:55 | 1 | UX80799.D | DB-624 0.18(mm) |
| STD8260 240-206828/9 IC | | 11/17/2015 06:16 | 1 | UX80800.D | DB-624 0.18(mm) |
| STD8260 240-206828/10 IC | | 11/17/2015 06:38 | 1 | UX80801.D | DB-624 0.18(mm) |
| ICV 240-206828/11 | | 11/17/2015 06:59 | 1 | UX80802.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/17/2015 07:41 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Instrument ID: A3UX8  Start Date: 11/20/2015 03:12

Analysis Batch Number: 207720  End Date: 11/20/2015 10:06

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-207720/1 | | 11/20/2015 03:12 | 1 | BFB8948.D | DB-624 0.18(mm) |
| CCVIS 240-207720/4 | | 11/20/2015 03:43 | 1 | UX80875.D | DB-624 0.18(mm) |
| CCV 240-207720/5 | | 11/20/2015 04:04 | 1 | UX80876.D | DB-624 0.18(mm) |
| LCS 240-207720/6 | | 11/20/2015 04:26 | 1 | UX80877.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/20/2015 04:47 | 1 | | DB-624 0.18(mm) |
| MB 240-207720/8 | | 11/20/2015 05:08 | 1 | UX80879.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/20/2015 06:55 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/20/2015 07:16 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/20/2015 07:37 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/20/2015 07:59 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/20/2015 08:20 | 1 | | DB-624 0.18(mm) |
| 240-57972-C-3-C MS | | 11/20/2015 08:41 | 1 | UX80889.D | DB-624 0.18(mm) |
| 240-57972-C-3-D MSD | | 11/20/2015 09:02 | 1 | UX80890.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/20/2015 09:24 | 1 | | DB-624 0.18(mm) |
| 240-57930-2 | F-VS-UST-P-2 | 11/20/2015 10:06 | 1 | UX80893.D | DB-624 0.18(mm) |

# GC/MS VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 207393  Batch Start Date: 11/18/15 12:27  Batch Analyst: Macenczak, Steven

Batch Method: 5030C  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57930-B-1 | F-VS-UST-P-1 | 5030C, 8260C | T | 5.18 g | 5 mL | | | | |
| 240-57930-B-3 | F-VS-UST-P-3 | 5030C, 8260C | T | 5.55 g | 5 mL | | | | |
| 240-57930-B-5 | F-VB-UST-P-1 | 5030C, 8260C | T | 5.06 g | 5 mL | | | | |
| 240-57930-B-6 | F-VB-UST-P-2 | 5030C, 8260C | T | 5.31 g | 5 mL | | | | |
| 240-57930-B-7 | F-VB-UST-P-3 | 5030C, 8260C | T | 5.29 g | 5 mL | | | | |
| 240-57930-B-8 | F-VB-UST-P-4 | 5030C, 8260C | T | 5.55 g | 5 mL | | | | |
| 240-57930-B-9 | F-VB-UST-P-5 | 5030C, 8260C | T | 5.55 g | 5 mL | | | | |
| 240-57930-B-10 | F-VB-UST-P-6 | 5030C, 8260C | T | 5.10 g | 5 mL | | | | |
| 240-57930-B-11 | F-VS-UST-P-DUP | 5030C, 8260C | T | 5.73 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| Balance ID | b050 |
| Blank Soil Lot Number | 148757 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC/MS VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 207613  Batch Start Date: 11/19/15 13:00  Batch Analyst: Macenczak, Steven

Batch Method: 5030C  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57930-B-2 | F-VS-UST-P-2 | 5030C, 8260C | T | 5.74 g | 5 mL | | | | |
| 240-57930-B-2 | F-VS-UST-P-2 | 5030C, 8260C | T | 5.09 g | 5 mL | | | | |
| 240-57930-B-4 | F-VS-UST-P-4 | 5030C, 8260C | T | 5.24 g | 5 mL | | | | |
| 240-57921-A-4-B MS | | 5030C, 8260C | T | 5.16 g | 5 mL | | | | |
| 240-57921-A-4-B MSD | | 5030C, 8260C | T | 5.24 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| | |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC/MS VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 207709  Batch Start Date: 11/20/15 04:52  Batch Analyst: Lata, Todd

Batch Method: 5030C  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57972-C-3 MS | | 5030C, 8260C | T | 5.28 g | 5 mL | | | | |
| 240-57972-C-3 MSD | | 5030C, 8260C | T | 5.04 g | 5 mL | | | | |
| 240-57930-B-2 | F-VS-UST-P-2 | 5030C, 8260C | T | 5.61 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| | |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8270D

## Semivolatile Organic Compounds (GC/MS) by Method 8270D

FORM II
GC/MS SEMI VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): DB-624 ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | 2FP # | PHL # | NBZ # | FBP # | TBP # | TPH # |
|---|---|---|---|---|---|---|---|
| F-VS-UST-P-1 | 240-57930-1 | 43 | 44 | 42 | 44 | 39 | 44 |
| F-VS-UST-P-2 | 240-57930-2 | 72 | 72 | 68 | 72 | 58 | 76 |
| F-VS-UST-P-3 | 240-57930-3 | 80 | 82 | 71 | 77 | 65 | 76 |
| F-VS-UST-P-4 | 240-57930-4 | 66 | 67 | 62 | 66 | 56 | 69 |
| F-VB-UST-P-1 | 240-57930-5 | 78 | 80 | 70 | 79 | 56 | 86 |
| F-VB-UST-P-2 | 240-57930-6 | 68 | 68 | 64 | 67 | 55 | 68 |
| F-VB-UST-P-3 | 240-57930-7 | 66 | 67 | 59 | 63 | 40 | 66 |
| F-VB-UST-P-4 | 240-57930-8 | 73 | 75 | 66 | 71 | 54 | 74 |
| F-VB-UST-P-5 | 240-57930-9 | 68 | 69 | 60 | 67 | 41 | 73 |
| F-VB-UST-P-6 | 240-57930-10 | 68 | 69 | 61 | 66 | 53 | 70 |
| F-VS-UST-P-DUP | 240-57930-11 | 68 | 68 | 65 | 66 | 62 | 66 |
| | MB 240-206887/23-A | 65 | 72 | 67 | 74 | 43 | 77 |
| | MB 240-206908/21-A | 70 | 73 | 67 | 70 | 30 | 74 |
| | LCS 240-206887/24-A | 80 | 80 | 73 | 77 | 69 | 78 |
| | LCS 240-206908/22-A | 70 | 71 | 70 | 70 | 42 | 71 |
| F-VB-UST-P-1 MS | 240-57930-5 MS | 78 | 78 | 73 | 74 | 72 | 76 |
| F-VB-UST-P-1 MSD | 240-57930-5 MSD | 73 | 74 | 71 | 72 | 64 | 75 |

| | QC LIMITS |
|---|---|
| 2FP = 2-Fluorophenol (Surr) | 24-110 |
| PHL = Phenol-d5 (Surr) | 26-110 |
| NBZ = Nitrobenzene-d5 (Surr) | 20-110 |
| FBP = 2-Fluorobiphenyl (Surr) | 24-110 |
| TBP = 2,4,6-Tribromophenol (Surr) | 10-110 |
| TPH = Terphenyl-d14 (Surr) | 36-110 |

# Column to be used to flag recovery values

FORM II 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51119008.D

Lab ID: LCS 240-206887/24-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 486 | 73 | 35-110 | |
| bis (2-chloroisopropyl) ether | 667 | 490 | 73 | 29-110 | |
| 2,4,5-Trichlorophenol | 667 | 484 | 73 | 25-110 | |
| 2,4,6-Trichlorophenol | 667 | 468 | 70 | 12-110 | |
| 2,4-Dichlorophenol | 667 | 499 | 75 | 39-110 | |
| 2,4-Dimethylphenol | 667 | 428 | 64 | 29-110 | |
| 2,4-Dinitrophenol | 1330 | 583 | 44 | 10-110 | |
| 2,4-Dinitrotoluene | 667 | 537 | 81 | 48-110 | |
| 2,6-Dinitrotoluene | 667 | 539 | 81 | 45-110 | |
| 2-Chloronaphthalene | 667 | 491 | 74 | 32-110 | |
| 2-Chlorophenol | 667 | 509 | 76 | 37-110 | |
| 2-Methylnaphthalene | 667 | 503 | 75 | 36-110 | |
| 2-Methylphenol | 667 | 500 | 75 | 41-110 | |
| 2-Nitroaniline | 667 | 507 | 76 | 45-110 | |
| 2-Nitrophenol | 667 | 509 | 76 | 34-110 | |
| 3,3'-Dichlorobenzidine | 1330 | 770 | 58 | 28-110 | |
| 3-Nitroaniline | 667 | 450 | 68 | 44-110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 727 | 55 | 10-110 | |
| 4-Bromophenyl phenyl ether | 667 | 539 | 81 | 39-110 | |
| 4-Chloro-3-methylphenol | 667 | 518 | 78 | 48-110 | |
| 4-Chloroaniline | 667 | 358 | 54 | 30-110 | |
| 4-Chlorophenyl phenyl ether | 667 | 513 | 77 | 40-110 | |
| 4-Nitroaniline | 667 | 534 | 80 | 48-110 | |
| 4-Nitrophenol | 1330 | 929 | 70 | 28-110 | |
| Acenaphthene | 667 | 487 | 73 | 38-110 | |
| Acenaphthylene | 667 | 498 | 75 | 40-110 | |
| Acetophenone | 667 | 491 | 74 | 40-110 | |
| Anthracene | 667 | 510 | 77 | 48-110 | |
| Atrazine | 1330 | 1060 | 79 | 54-120 | |
| Benzaldehyde | 1330 | 1010 | 76 | 32-110 | |
| Benzo[a]anthracene | 667 | 517 | 78 | 50-110 | |
| Benzo[a]pyrene | 667 | 509 | 76 | 44-110 | |
| Benzo[b]fluoranthene | 667 | 509 | 76 | 43-110 | |
| Benzo[g,h,i]perylene | 667 | 518 | 78 | 51-110 | |
| Benzo[k]fluoranthene | 667 | 530 | 79 | 38-105 | |
| Bis(2-chloroethoxy)methane | 667 | 526 | 79 | 32-110 | |
| Bis(2-chloroethyl)ether | 667 | 515 | 77 | 34-110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 607 | 91 | 50-110 | |
| Butyl benzyl phthalate | 667 | 545 | 82 | 51-110 | |
| Caprolactam | 1330 | 1100 | 82 | 44-114 | |
| Carbazole | 667 | 530 | 80 | 50-110 | |
| Chrysene | 667 | 538 | 81 | 50-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51119008.D

Lab ID: LCS 240-206887/24-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 667 | 504 | 76 | 51-110 | |
| Dibenzofuran | 667 | 499 | 75 | 43-110 | |
| Diethyl phthalate | 667 | 544 | 82 | 52-110 | |
| Dimethyl phthalate | 667 | 523 | 79 | 50-110 | |
| Di-n-butyl phthalate | 667 | 549 | 82 | 51-110 | |
| Di-n-octyl phthalate | 667 | 515 | 77 | 48-110 | |
| Fluoranthene | 667 | 528 | 79 | 51-110 | |
| Fluorene | 667 | 491 | 74 | 46-110 | |
| Hexachlorobenzene | 667 | 518 | 78 | 43-110 | |
| Hexachlorobutadiene | 667 | 490 | 73 | 29-110 | |
| Hexachlorocyclopentadiene | 667 | 320 J | 48 | 12-110 | |
| Hexachloroethane | 667 | 473 | 71 | 30-110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 515 | 77 | 50-110 | |
| Isophorone | 667 | 527 | 79 | 36-110 | |
| N-Nitrosodi-n-propylamine | 667 | 513 | 77 | 38-110 | |
| N-Nitrosodiphenylamine | 1330 | 1030 | 77 | 46-110 | |
| Naphthalene | 667 | 505 | 76 | 36-110 | |
| Nitrobenzene | 667 | 508 | 76 | 32-110 | |
| Pentachlorophenol | 1330 | 665 | 50 | 10-110 | |
| Phenanthrene | 667 | 516 | 77 | 49-110 | |
| Phenol | 667 | 509 | 76 | 38-110 | |
| Pyrene | 667 | 507 | 76 | 49-110 | |
| 3 & 4 Methylphenol | 667 | 501 | 75 | 40-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51118004.D

Lab ID: LCS 240-206908/22-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 442 | 66 | 35-110 | |
| bis (2-chloroisopropyl) ether | 667 | 446 | 67 | 29-110 | |
| 2,4,5-Trichlorophenol | 667 | 387 | 58 | 25-110 | |
| 2,4,6-Trichlorophenol | 667 | 286 | 43 | 12-110 | |
| 2,4-Dichlorophenol | 667 | 461 | 69 | 39-110 | |
| 2,4-Dimethylphenol | 667 | 467 | 70 | 29-110 | |
| 2,4-Dinitrophenol | 1330 | 611 | 46 | 10-110 | |
| 2,4-Dinitrotoluene | 667 | 488 | 73 | 48-110 | |
| 2,6-Dinitrotoluene | 667 | 480 | 72 | 45-110 | |
| 2-Chloronaphthalene | 667 | 463 | 69 | 32-110 | |
| 2-Chlorophenol | 667 | 445 | 67 | 37-110 | |
| 2-Methylnaphthalene | 667 | 458 | 69 | 36-110 | |
| 2-Methylphenol | 667 | 453 | 68 | 41-110 | |
| 2-Nitroaniline | 667 | 485 | 73 | 45-110 | |
| 2-Nitrophenol | 667 | 461 | 69 | 34-110 | |
| 3,3'-Dichlorobenzidine | 1330 | 732 | 55 | 28-110 | |
| 3-Nitroaniline | 667 | 432 | 65 | 44-110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 779 | 58 | 10-110 | |
| 4-Bromophenyl phenyl ether | 667 | 488 | 73 | 39-110 | |
| 4-Chloro-3-methylphenol | 667 | 474 | 71 | 48-110 | |
| 4-Chloroaniline | 667 | 349 | 52 | 30-110 | |
| 4-Chlorophenyl phenyl ether | 667 | 467 | 70 | 40-110 | |
| 4-Nitroaniline | 667 | 503 | 75 | 48-110 | |
| 4-Nitrophenol | 1330 | 881 | 66 | 28-110 | |
| Acenaphthene | 667 | 464 | 70 | 38-110 | |
| Acenaphthylene | 667 | 462 | 69 | 40-110 | |
| Acetophenone | 667 | 438 | 66 | 40-110 | |
| Anthracene | 667 | 471 | 71 | 48-110 | |
| Atrazine | 1330 | 971 | 73 | 54-120 | |
| Benzaldehyde | 1330 | 909 | 68 | 32-110 | |
| Benzo[a]anthracene | 667 | 466 | 70 | 50-110 | |
| Benzo[a]pyrene | 667 | 466 | 70 | 44-110 | |
| Benzo[b]fluoranthene | 667 | 452 | 68 | 43-110 | |
| Benzo[g,h,i]perylene | 667 | 476 | 71 | 51-110 | |
| Benzo[k]fluoranthene | 667 | 469 | 70 | 38-105 | |
| Bis(2-chloroethoxy)methane | 667 | 486 | 73 | 32-110 | |
| Bis(2-chloroethyl)ether | 667 | 453 | 68 | 34-110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 704 | 106 | 50-110 | |
| Butyl benzyl phthalate | 667 | 489 | 73 | 51-110 | |
| Caprolactam | 1330 | 1050 | 79 | 44-114 | |
| Carbazole | 667 | 488 | 73 | 50-110 | |
| Chrysene | 667 | 495 | 74 | 50-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51118004.D

Lab ID: LCS 240-206908/22-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 667 | 469 | 70 | 51-110 | |
| Dibenzofuran | 667 | 464 | 70 | 43-110 | |
| Diethyl phthalate | 667 | 490 | 73 | 52-110 | |
| Dimethyl phthalate | 667 | 479 | 72 | 50-110 | |
| Di-n-butyl phthalate | 667 | 489 | 73 | 51-110 | |
| Di-n-octyl phthalate | 667 | 445 | 67 | 48-110 | |
| Fluoranthene | 667 | 473 | 71 | 51-110 | |
| Fluorene | 667 | 454 | 68 | 46-110 | |
| Hexachlorobenzene | 667 | 477 | 72 | 43-110 | |
| Hexachlorobutadiene | 667 | 447 | 67 | 29-110 | |
| Hexachlorocyclopentadiene | 667 | 397 | 60 | 12-110 | |
| Hexachloroethane | 667 | 418 | 63 | 30-110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 477 | 72 | 50-110 | |
| Isophorone | 667 | 478 | 72 | 36-110 | |
| N-Nitrosodi-n-propylamine | 667 | 451 | 68 | 38-110 | |
| N-Nitrosodiphenylamine | 1330 | 946 | 71 | 46-110 | |
| Naphthalene | 667 | 462 | 69 | 36-110 | |
| Nitrobenzene | 667 | 467 | 70 | 32-110 | |
| Pentachlorophenol | 1330 | 669 | 50 | 10-110 | |
| Phenanthrene | 667 | 471 | 71 | 49-110 | |
| Phenol | 667 | 456 | 68 | 38-110 | |
| Pyrene | 667 | 475 | 71 | 49-110 | |
| 3 & 4 Methylphenol | 667 | 448 | 67 | 40-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51119015.D

Lab ID: 240-57930-5 MS Client ID: F-VB-UST-P-1 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 786 | 4.1 U | 547 | 70 | 32-110 | |
| bis (2-chloroisopropyl) ether | 786 | 11 U | 561 | 71 | 11-110 | |
| 2,4,5-Trichlorophenol | 786 | 29 U | 554 | 70 | 10-117 | |
| 2,4,6-Trichlorophenol | 786 | 10 U | 543 | 69 | 10-110 | |
| 2,4-Dichlorophenol | 786 | 23 U | 567 | 72 | 10-110 | |
| 2,4-Dimethylphenol | 786 | 23 U | 554 | 71 | 10-110 | |
| 2,4-Dinitrophenol | 1570 | 25 U | 752 | 48 | 10-110 | |
| 2,4-Dinitrotoluene | 786 | 20 U | 582 | 74 | 32-110 | |
| 2,6-Dinitrotoluene | 786 | 25 U | 567 | 72 | 35-110 | |
| 2-Chloronaphthalene | 786 | 0.53 U | 547 | 70 | 28-110 | |
| 2-Chlorophenol | 786 | 9.6 U | 587 | 75 | 10-110 | |
| 2-Methylnaphthalene | 786 | 0.59 U | 564 | 72 | 10-133 | |
| 2-Methylphenol | 786 | 13 U | 587 | 75 | 24-110 | |
| 2-Nitroaniline | 786 | 11 U | 558 | 71 | 39-110 | |
| 2-Nitrophenol | 786 | 9.7 U | 553 | 70 | 10-110 | |
| 3,3'-Dichlorobenzidine | 1570 | 21 U | 856 | 54 | 10-110 | |
| 3-Nitroaniline | 786 | 19 U | 491 | 63 | 10-110 | |
| 4,6-Dinitro-2-methylphenol | 1570 | 11 U | 856 | 54 | 10-110 | |
| 4-Bromophenyl phenyl ether | 786 | 15 U | 597 | 76 | 33-110 | |
| 4-Chloro-3-methylphenol | 786 | 25 U | 598 | 76 | 25-110 | |
| 4-Chloroaniline | 786 | 20 U | 401 | 51 | 10-110 | |
| 4-Chlorophenyl phenyl ether | 786 | 15 U | 567 | 72 | 32-110 | |
| 4-Nitroaniline | 786 | 30 U | 510 | 65 | 10-110 | |
| 4-Nitrophenol | 1570 | 20 U | 1030 | 65 | 10-113 | |
| Acenaphthene | 786 | 0.89 U | 557 | 71 | 22-110 | |
| Acenaphthylene | 786 | 0.41 U | 562 | 71 | 24-110 | |
| Acetophenone | 786 | 11 U | 559 | 71 | 31-110 | |
| Anthracene | 786 | 0.91 U | 552 | 70 | 20-110 | |
| Atrazine | 1570 | 11 U | 1150 | 73 | 45-118 | |
| Benzaldehyde | 1570 | 14 U | 1160 | 74 | 23-110 | |
| Benzo[a]anthracene | 786 | 0.74 U | 567 | 72 | 10-122 | |
| Benzo[a]pyrene | 786 | 0.75 U | 564 | 72 | 10-110 | |
| Benzo[b]fluoranthene | 786 | 0.69 U | 560 | 71 | 12-118 | |
| Benzo[g,h,i]perylene | 786 | 0.41 U | 561 | 71 | 10-117 | |
| Benzo[k]fluoranthene | 786 | 0.80 U | 576 | 73 | 10-121 | |
| Bis(2-chloroethoxy)methane | 786 | 26 U | 605 | 77 | 26-110 | |
| Bis(2-chloroethyl)ether | 786 | 2.3 U | 593 | 75 | 21-110 | |
| Bis(2-ethylhexyl) phthalate | 786 | 25 J | 613 | 75 | 40-110 | |
| Butyl benzyl phthalate | 786 | 12 U | 601 | 77 | 44-110 | |
| Caprolactam | 1570 | 43 U | 1300 | 83 | 10-134 | |
| Carbazole | 786 | 32 U | 537 | 68 | 34-110 | |
| Chrysene | 786 | 1.3 U | 586 | 75 | 10-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51119015.D

Lab ID: 240-57930-5 MS Client ID: F-VB-UST-P-1 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 786 | 0.77 U | 564 | 72 | 14-113 | |
| Dibenzofuran | 786 | 0.77 U | 562 | 71 | 29-110 | |
| Diethyl phthalate | 786 | 19 U | 591 | 75 | 42-110 | |
| Dimethyl phthalate | 786 | 20 U | 575 | 73 | 41-110 | |
| Di-n-butyl phthalate | 786 | 18 U | 585 | 74 | 43-110 | |
| Di-n-octyl phthalate | 786 | 9.3 U | 570 | 73 | 24-119 | |
| Fluoranthene | 786 | 0.64 U | 574 | 73 | 10-110 | |
| Fluorene | 786 | 0.62 U | 553 | 70 | 23-110 | |
| Hexachlorobenzene | 786 | 2.5 U | 551 | 70 | 34-110 | |
| Hexachlorobutadiene | 786 | 6.6 U | 543 | 69 | 25-110 | |
| Hexachlorocyclopentadiene | 786 | 9.5 U | 315 J | 40 | 10-110 | |
| Hexachloroethane | 786 | 11 U | 519 | 66 | 12-110 | |
| Indeno[1,2,3-cd]pyrene | 786 | 0.41 U | 558 | 71 | 10-114 | |
| Isophorone | 786 | 15 U | 596 | 76 | 29-110 | |
| N-Nitrosodi-n-propylamine | 786 | 7.4 U | 570 | 72 | 26-110 | |
| N-Nitrosodiphenylamine | 1570 | 25 U | 1140 | 72 | 22-110 | |
| Naphthalene | 786 | 0.96 U | 563 | 72 | 10-111 | |
| Nitrobenzene | 786 | 2.6 U | 571 | 73 | 23-110 | |
| Pentachlorophenol | 1570 | 11 U | 863 | 55 | 10-110 | |
| Phenanthrene | 786 | 0.86 U | 560 | 71 | 10-166 | |
| Phenol | 786 | 8.6 U | 583 | 74 | 17-110 | |
| Pyrene | 786 | 0.52 U | 579 | 74 | 10-147 | |
| 3 & 4 Methylphenol | 786 | 23 U | 578 | 74 | 25-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51119016.D

Lab ID: 240-57930-5 MSD Client ID: F-VB-UST-P-1 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 791 | 534 | 67 | 3 | 32 | 32-110 | |
| bis (2-chloroisopropyl) ether | 791 | 534 | 68 | 5 | 42 | 11-110 | |
| 2,4,5-Trichlorophenol | 791 | 553 | 70 | 0 | 99 | 10-117 | |
| 2,4,6-Trichlorophenol | 791 | 506 | 64 | 7 | 38 | 10-110 | |
| 2,4-Dichlorophenol | 791 | 568 | 72 | 0 | 34 | 10-110 | |
| 2,4-Dimethylphenol | 791 | 535 | 68 | 4 | 31 | 10-110 | |
| 2,4-Dinitrophenol | 1580 | 741 | 47 | 1 | 99 | 10-110 | |
| 2,4-Dinitrotoluene | 791 | 593 | 75 | 2 | 30 | 32-110 | |
| 2,6-Dinitrotoluene | 791 | 567 | 72 | 0 | 30 | 35-110 | |
| 2-Chloronaphthalene | 791 | 545 | 69 | 0 | 30 | 28-110 | |
| 2-Chlorophenol | 791 | 546 | 69 | 7 | 47 | 10-110 | |
| 2-Methylnaphthalene | 791 | 568 | 72 | 1 | 42 | 10-133 | |
| 2-Methylphenol | 791 | 554 | 70 | 6 | 51 | 24-110 | |
| 2-Nitroaniline | 791 | 568 | 72 | 2 | 31 | 39-110 | |
| 2-Nitrophenol | 791 | 555 | 70 | 0 | 49 | 10-110 | |
| 3,3'-Dichlorobenzidine | 1580 | 899 | 57 | 5 | 56 | 10-110 | |
| 3-Nitroaniline | 791 | 483 | 61 | 2 | 30 | 10-110 | |
| 4,6-Dinitro-2-methylphenol | 1580 | 875 | 55 | 2 | 55 | 10-110 | |
| 4-Bromophenyl phenyl ether | 791 | 593 | 75 | 1 | 30 | 33-110 | |
| 4-Chloro-3-methylphenol | 791 | 579 | 73 | 3 | 54 | 25-110 | |
| 4-Chloroaniline | 791 | 419 | 53 | 4 | 36 | 10-110 | |
| 4-Chlorophenyl phenyl ether | 791 | 571 | 72 | 1 | 30 | 32-110 | |
| 4-Nitroaniline | 791 | 524 | 66 | 3 | 48 | 10-110 | |
| 4-Nitrophenol | 1580 | 1020 | 64 | 1 | 49 | 10-113 | |
| Acenaphthene | 791 | 542 | 69 | 3 | 99 | 22-110 | |
| Acenaphthylene | 791 | 552 | 70 | 2 | 99 | 24-110 | |
| Acetophenone | 791 | 538 | 68 | 4 | 43 | 31-110 | |
| Anthracene | 791 | 564 | 71 | 2 | 99 | 20-110 | |
| Atrazine | 1580 | 1190 | 75 | 4 | 30 | 45-118 | |
| Benzaldehyde | 1580 | 1100 | 69 | 5 | 42 | 23-110 | |
| Benzo[a]anthracene | 791 | 580 | 73 | 2 | 99 | 10-122 | |
| Benzo[a]pyrene | 791 | 588 | 74 | 4 | 99 | 10-110 | |
| Benzo[b]fluoranthene | 791 | 575 | 73 | 3 | 99 | 12-118 | |
| Benzo[g,h,i]perylene | 791 | 571 | 72 | 2 | 99 | 10-117 | |
| Benzo[k]fluoranthene | 791 | 598 | 76 | 4 | 99 | 10-121 | |
| Bis(2-chloroethoxy)methane | 791 | 594 | 75 | 2 | 37 | 26-110 | |
| Bis(2-chloroethyl)ether | 791 | 564 | 71 | 5 | 55 | 21-110 | |
| Bis(2-ethylhexyl) phthalate | 791 | 627 | 76 | 2 | 30 | 40-110 | |
| Butyl benzyl phthalate | 791 | 617 | 78 | 3 | 30 | 44-110 | |
| Caprolactam | 1580 | 1320 | 83 | 1 | 32 | 10-134 | |
| Carbazole | 791 | 545 | 69 | 1 | 30 | 34-110 | |
| Chrysene | 791 | 608 | 77 | 4 | 99 | 10-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 51119016.D

Lab ID: 240-57930-5 MSD  Client ID: F-VB-UST-P-1 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 791 | 575 | 73 | 2 | 99 | 14-113 | |
| Dibenzofuran | 791 | 552 | 70 | 2 | 30 | 29-110 | |
| Diethyl phthalate | 791 | 590 | 75 | 0 | 30 | 42-110 | |
| Dimethyl phthalate | 791 | 576 | 73 | 0 | 30 | 41-110 | |
| Di-n-butyl phthalate | 791 | 593 | 75 | 1 | 30 | 43-110 | |
| Di-n-octyl phthalate | 791 | 592 | 75 | 4 | 30 | 24-119 | |
| Fluoranthene | 791 | 578 | 73 | 1 | 99 | 10-110 | |
| Fluorene | 791 | 545 | 69 | 1 | 99 | 23-110 | |
| Hexachlorobenzene | 791 | 562 | 71 | 2 | 30 | 34-110 | |
| Hexachlorobutadiene | 791 | 548 | 69 | 1 | 34 | 25-110 | |
| Hexachlorocyclopentadiene | 791 | 238 J | 30 | 28 | 79 | 10-110 | |
| Hexachloroethane | 791 | 485 | 61 | 7 | 50 | 12-110 | |
| Indeno[1,2,3-cd]pyrene | 791 | 579 | 73 | 4 | 99 | 10-114 | |
| Isophorone | 791 | 586 | 74 | 2 | 38 | 29-110 | |
| N-Nitrosodi-n-propylamine | 791 | 540 | 68 | 5 | 42 | 26-110 | |
| N-Nitrosodiphenylamine | 1580 | 1130 | 71 | 1 | 30 | 22-110 | |
| Naphthalene | 791 | 559 | 71 | 1 | 99 | 10-111 | |
| Nitrobenzene | 791 | 562 | 71 | 2 | 41 | 23-110 | |
| Pentachlorophenol | 1580 | 793 | 50 | 8 | 50 | 10-110 | |
| Phenanthrene | 791 | 567 | 72 | 1 | 99 | 10-166 | |
| Phenol | 791 | 554 | 70 | 5 | 53 | 17-110 | |
| Pyrene | 791 | 586 | 74 | 1 | 99 | 10-147 | |
| 3 & 4 Methylphenol | 791 | 544 | 69 | 6 | 50 | 25-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM IV
GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: 51119007.D  Lab Sample ID: MB 240-206887/23-A

Matrix: Solid  Date Extracted: 11/16/2015 08:48

Instrument ID: A4HP10  Date Analyzed: 11/19/2015 11:50

Level:(Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-206887/24-A | 51119008.D | 11/19/2015 12:13 |
| F-VB-UST-P-2 | 240-57930-6 | 51119017.D | 11/19/2015 15:48 |
| F-VB-UST-P-3 | 240-57930-7 | 51119018.D | 11/19/2015 16:12 |
| F-VB-UST-P-4 | 240-57930-8 | 51119019.D | 11/19/2015 16:36 |
| F-VB-UST-P-5 | 240-57930-9 | 51119020.D | 11/19/2015 17:00 |
| F-VB-UST-P-6 | 240-57930-10 | 51119021.D | 11/19/2015 17:23 |
| F-VS-UST-P-DUP | 240-57930-11 | 51119028.D | 11/19/2015 20:10 |

FORM IV
GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: 51118003.D  Lab Sample ID: MB 240-206908/21-A

Matrix: Solid  Date Extracted: 11/16/2015 09:20

Instrument ID: A4HP10  Date Analyzed: 11/18/2015 08:51

Level:(Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-206908/22-A | 51118004.D | 11/18/2015 09:15 |
| F-VB-UST-P-1 | 240-57930-5 | 51119014.D | 11/19/2015 14:37 |
| F-VB-UST-P-1 MS | 240-57930-5 MS | 51119015.D | 11/19/2015 15:00 |
| F-VB-UST-P-1 MSD | 240-57930-5 MSD | 51119016.D | 11/19/2015 15:24 |
| F-VS-UST-P-2 | 240-57930-2 | 51119026.D | 11/19/2015 19:23 |
| F-VS-UST-P-1 | 240-57930-1 | 51119027.D | 11/19/2015 19:46 |
| F-VS-UST-P-4 | 240-57930-4 | 51119029.D | 11/19/2015 20:34 |
| F-VS-UST-P-3 | 240-57930-3 | 51119030.D | 11/19/2015 20:57 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: 51113001.D  DFTPP Injection Date: 11/13/2015

Instrument ID: A4HP10  DFTPP Injection Time: 09:10

Analysis Batch No.: 206629

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 46.1 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0) 1 |
| 69 | Mass 69 relative abundance | 37.7 | |
| 70 | Less than 2.0% of mass 69 | 0.2 | (0.5) 1 |
| 127 | 25.0 - 75.0% of mass 198 | 52.9 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.8 | |
| 275 | 10.0- 30.0% of mass 198 | 24.8 | |
| 365 | Greater than 0.75% of mass 198 | 3.4 | |
| 441 | Present, but less than mass 443 | 15.1 | |
| 442 | 40.0 - 110.0% of mass 198 | 92.2 | |
| 443 | 15.0 - 24.0% of mass 442 | 18.4 | (20.0) 2 |

1-Value is % mass 69  2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD5 240-206629/2 | 51113002.D | 11/13/2015 | 09:35 |
| | STD4 240-206629/3 | 51113003.D | 11/13/2015 | 09:59 |
| | STD3 240-206629/4 | 51113004.D | 11/13/2015 | 10:23 |
| | STD2 240-206629/5 | 51113005.D | 11/13/2015 | 10:48 |
| | STD1 240-206629/6 | 51113006.D | 11/13/2015 | 11:12 |
| | STD8 240-206629/8 | 51113008.D | 11/13/2015 | 12:00 |
| | STD7 240-206629/9 | 51113009.D | 11/13/2015 | 12:24 |
| | STD6 240-206629/10 | 51113010.D | 11/13/2015 | 12:48 |
| | ICV 240-206629/11 | 51113011.D | 11/13/2015 | 13:12 |
| | STD9 240-206629/12 | 51113012.D | 11/13/2015 | 13:36 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: 51118001.D  DFTPP Injection Date: 11/18/2015

Instrument ID: A4HP10  DFTPP Injection Time: 08:03

Analysis Batch No.: 207287

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 44.6 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0) 1 |
| 69 | Mass 69 relative abundance | 37.3 | |
| 70 | Less than 2.0% of mass 69 | 0.2 | (0.5) 1 |
| 127 | 25.0 - 75.0% of mass 198 | 52.0 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.6 | |
| 275 | 10.0- 30.0% of mass 198 | 24.6 | |
| 365 | Greater than 0.75% of mass 198 | 3.4 | |
| 441 | Present, but less than mass 443 | 15.8 | |
| 442 | 40.0 - 110.0% of mass 198 | 96.1 | |
| 443 | 15.0 - 24.0% of mass 442 | 19.1 | (19.9) 2 |

1-Value is % mass 69  2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCV 240-207287/2 | 51118002.D | 11/18/2015 | 08:26 |
| | MB 240-206908/21-A | 51118003.D | 11/18/2015 | 08:51 |
| | LCS 240-206908/22-A | 51118004.D | 11/18/2015 | 09:15 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab File ID: 51119001.D  DFTPP Injection Date: 11/19/2015

Instrument ID: A4HP10  DFTPP Injection Time: 09:29

Analysis Batch No.: 207541

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 42.9 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0) 1 |
| 69 | Mass 69 relative abundance | 36.7 | |
| 70 | Less than 2.0% of mass 69 | 0.2 | (0.5) 1 |
| 127 | 25.0 - 75.0% of mass 198 | 52.5 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.7 | |
| 275 | 10.0- 30.0% of mass 198 | 25.0 | |
| 365 | Greater than 0.75% of mass 198 | 3.6 | |
| 441 | Present, but less than mass 443 | 17.1 | |
| 442 | 40.0 - 110.0% of mass 198 | 100.0 | |
| 443 | 15.0 - 24.0% of mass 442 | 19.5 | (19.5) 2 |

1-Value is % mass 69  2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCV 240-207541/2 | 51119002.D | 11/19/2015 | 09:50 |
| | MB 240-206887/23-A | 51119007.D | 11/19/2015 | 11:50 |
| | LCS 240-206887/24-A | 51119008.D | 11/19/2015 | 12:13 |
| F-VB-UST-P-1 | 240-57930-5 | 51119014.D | 11/19/2015 | 14:37 |
| F-VB-UST-P-1 MS | 240-57930-5 MS | 51119015.D | 11/19/2015 | 15:00 |
| F-VB-UST-P-1 MSD | 240-57930-5 MSD | 51119016.D | 11/19/2015 | 15:24 |
| F-VB-UST-P-2 | 240-57930-6 | 51119017.D | 11/19/2015 | 15:48 |
| F-VB-UST-P-3 | 240-57930-7 | 51119018.D | 11/19/2015 | 16:12 |
| F-VB-UST-P-4 | 240-57930-8 | 51119019.D | 11/19/2015 | 16:36 |
| F-VB-UST-P-5 | 240-57930-9 | 51119020.D | 11/19/2015 | 17:00 |
| F-VB-UST-P-6 | 240-57930-10 | 51119021.D | 11/19/2015 | 17:23 |
| F-VS-UST-P-2 | 240-57930-2 | 51119026.D | 11/19/2015 | 19:23 |
| F-VS-UST-P-1 | 240-57930-1 | 51119027.D | 11/19/2015 | 19:46 |
| F-VS-UST-P-DUP | 240-57930-11 | 51119028.D | 11/19/2015 | 20:10 |
| F-VS-UST-P-4 | 240-57930-4 | 51119029.D | 11/19/2015 | 20:34 |
| F-VS-UST-P-3 | 240-57930-3 | 51119030.D | 11/19/2015 | 20:57 |

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: STD6 240-206629/10  Date Analyzed: 11/13/2015 12:48

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 51113010.D  Heated Purge: (Y/N) N

Calibration ID: 31440

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 167224 | 6.09 | 697175 | 7.21 | 407101 | 8.71 |
| UPPER LIMIT | | 334448 | 6.59 | 1394350 | 7.71 | 814202 | 9.21 |
| LOWER LIMIT | | 83612 | 5.59 | 348588 | 6.71 | 203551 | 8.21 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-206629/11 | | 190384 | 6.09 | 768772 | 7.21 | 440245 | 8.71 |
| CCV 240-207287/2 CCVIS | | 159115 | 6.07 | 630914 | 7.19 | 366314 | 8.69 |
| CCV 240-207541/2 CCVIS | | 161328 | 6.06 | 665148 | 7.18 | 388586 | 8.68 |

DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Sample No.: STD6 240-206629/10 Date Analyzed: 11/13/2015 12:48

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm)

Lab File ID (Standard): 51113010.D Heated Purge: (Y/N) N

Calibration ID: 31440

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 626897 | 9.98 | 718468 | 12.61 | 611718 | 14.75 |
| UPPER LIMIT | | 1253794 | 10.48 | 1436936 | 13.11 | 1223436 | 15.25 |
| LOWER LIMIT | | 313449 | 9.48 | 359234 | 12.11 | 305859 | 14.25 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-206629/11 | | 672676 | 9.98 | 777993 | 12.60 | 638595 | 14.75 |
| CCV 240-207287/2 CCVIS | | 561626 | 9.95 | 660749 | 12.57 | 555821 | 14.70 |
| CCV 240-207541/2 CCVIS | | 619847 | 9.94 | 722332 | 12.55 | 599867 | 14.67 |

PHN = Phenanthrene-d10
CRY = Chrysene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Sample No.: CCV 240-207287/2 Date Analyzed: 11/18/2015 08:26

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm)

Lab File ID (Standard): 51118002.D Heated Purge: (Y/N) N

Calibration ID: 31440

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 159115 | 6.07 | 630914 | 7.19 | 366314 | 8.69 |
| UPPER LIMIT | | 318230 | 6.57 | 1261828 | 7.69 | 732628 | 9.19 |
| LOWER LIMIT | | 79558 | 5.57 | 315457 | 6.69 | 183157 | 8.19 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-206908/21-A | | 165580 | 6.08 | 667132 | 7.19 | 403782 | 8.69 |
| LCS 240-206908/22-A | | 156476 | 6.08 | 616302 | 7.19 | 362629 | 8.69 |

DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCV 240-207287/2  Date Analyzed: 11/18/2015 08:26

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 51118002.D  Heated Purge: (Y/N) N

Calibration ID: 31440

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 561626 | 9.95 | 660749 | 12.57 | 555821 | 14.70 |
| UPPER LIMIT | | 1123252 | 10.45 | 1321498 | 13.07 | 1111642 | 15.20 |
| LOWER LIMIT | | 280813 | 9.45 | 330375 | 12.07 | 277911 | 14.20 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-206908/21-A | | 626122 | 9.95 | 709483 | 12.56 | 631387 | 14.70 |
| LCS 240-206908/22-A | | 569523 | 9.95 | 655734 | 12.57 | 587170 | 14.70 |

PHN = Phenanthrene-d10
CRY = Chrysene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCV 240-207541/2  Date Analyzed: 11/19/2015 09:50

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 51119002.D  Heated Purge: (Y/N) N

Calibration ID: 31440

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 161328 | 6.06 | 665148 | 7.18 | 388586 | 8.68 |
| UPPER LIMIT | | 322656 | 6.56 | 1330296 | 7.68 | 777172 | 9.18 |
| LOWER LIMIT | | 80664 | 5.56 | 332574 | 6.68 | 194293 | 8.18 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-206887/23-A | | 147780 | 6.06 | 619471 | 7.18 | 373431 | 8.67 |
| LCS 240-206887/24-A | | 149366 | 6.06 | 615257 | 7.18 | 376589 | 8.67 |
| 240-57930-5 | F-VB-UST-P-1 | 153611 | 6.06 | 619772 | 7.18 | 369895 | 8.67 |
| 240-57930-5 MS | F-VB-UST-P-1 MS | 153372 | 6.06 | 637619 | 7.18 | 388418 | 8.67 |
| 240-57930-5 MSD | F-VB-UST-P-1 MSD | 164720 | 6.06 | 658981 | 7.18 | 405255 | 8.67 |
| 240-57930-6 | F-VB-UST-P-2 | 149551 | 6.06 | 598332 | 7.18 | 367668 | 8.67 |
| 240-57930-7 | F-VB-UST-P-3 | 151252 | 6.06 | 629109 | 7.18 | 381483 | 8.67 |
| 240-57930-8 | F-VB-UST-P-4 | 148845 | 6.06 | 621963 | 7.18 | 382651 | 8.67 |
| 240-57930-9 | F-VB-UST-P-5 | 151439 | 6.06 | 625362 | 7.18 | 387308 | 8.67 |
| 240-57930-10 | F-VB-UST-P-6 | 155865 | 6.06 | 642935 | 7.18 | 396330 | 8.67 |
| 240-57930-2 | F-VS-UST-P-2 | 163482 | 6.06 | 636886 | 7.18 | 384461 | 8.67 |
| 240-57930-1 | F-VS-UST-P-1 | 164323 | 6.06 | 659617 | 7.18 | 409587 | 8.67 |
| 240-57930-11 | F-VS-UST-P-DUP | 160354 | 6.06 | 645673 | 7.18 | 393379 | 8.67 |
| 240-57930-4 | F-VS-UST-P-4 | 159533 | 6.06 | 637723 | 7.18 | 394299 | 8.67 |
| 240-57930-3 | F-VS-UST-P-3 | 149741 | 6.06 | 630088 | 7.18 | 388571 | 8.67 |

DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCV 240-207541/2  Date Analyzed: 11/19/2015 09:50

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 51119002.D  Heated Purge: (Y/N) N

Calibration ID: 31440

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 619847 | 9.94 | 722332 | 12.55 | 599867 | 14.67 |
| UPPER LIMIT | | 1239694 | 10.44 | 1444664 | 13.05 | 1199734 | 15.17 |
| LOWER LIMIT | | 309924 | 9.44 | 361166 | 12.05 | 299934 | 14.17 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-206887/23-A | | 590416 | 9.94 | 666761 | 12.55 | 564893 | 14.67 |
| LCS 240-206887/24-A | | 593056 | 9.94 | 695958 | 12.55 | 601622 | 14.67 |
| 240-57930-5 | F-VB-UST-P-1 | 587594 | 9.94 | 656454 | 12.55 | 428238 | 14.67 |
| 240-57930-5 MS | F-VB-UST-P-1 MS | 611945 | 9.94 | 703523 | 12.55 | 612174 | 14.67 |
| 240-57930-5 MSD | F-VB-UST-P-1 MSD | 638125 | 9.94 | 734923 | 12.55 | 625674 | 14.67 |
| 240-57930-6 | F-VB-UST-P-2 | 581450 | 9.94 | 648887 | 12.55 | 506456 | 14.67 |
| 240-57930-7 | F-VB-UST-P-3 | 588779 | 9.94 | 667351 | 12.55 | 513166 | 14.67 |
| 240-57930-8 | F-VB-UST-P-4 | 590966 | 9.94 | 668250 | 12.55 | 538037 | 14.67 |
| 240-57930-9 | F-VB-UST-P-5 | 579702 | 9.94 | 664684 | 12.55 | 524803 | 14.67 |
| 240-57930-10 | F-VB-UST-P-6 | 614861 | 9.94 | 689315 | 12.55 | 546172 | 14.67 |
| 240-57930-2 | F-VS-UST-P-2 | 605473 | 9.94 | 685391 | 12.55 | 566491 | 14.67 |
| 240-57930-1 | F-VS-UST-P-1 | 656345 | 9.94 | 730391 | 12.55 | 580328 | 14.67 |
| 240-57930-11 | F-VS-UST-P-DUP | 625198 | 9.94 | 719073 | 12.55 | 536626 | 14.68 |
| 240-57930-4 | F-VS-UST-P-4 | 622150 | 9.94 | 709030 | 12.55 | 522377 | 14.68 |
| 240-57930-3 | F-VS-UST-P-3 | 626391 | 9.94 | 716801 | 12.55 | 536977 | 14.68 |

PHN = Phenanthrene-d10
CRY = Chrysene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-1 | Lab Sample ID: 240-57930-1

Matrix: Solid | Lab File ID: 51119027.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:30

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 29.52(g) | Date Analyzed: 11/19/2015 19:46

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 60 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.54 | U | 60 | 0.54 |
| 95-57-8 | 2-Chlorophenol | 9.8 | U | 60 | 9.8 |
| 91-57-6 | 2-Methylnaphthalene | 7.1 | J | 8.0 | 0.60 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.9 | U | 60 | 9.9 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 60 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 60 | 16 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.91 | U | 8.0 | 0.91 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 8.0 | 0.42 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 4.6 | J | 8.0 | 0.93 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 17 | J | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 12 | | 8.0 | 0.75 |
| 50-32-8 | Benzo[a]pyrene | 12 | | 8.0 | 0.77 |
| 205-99-2 | Benzo[b]fluoranthene | 17 | | 8.0 | 0.71 |
| 191-24-2 | Benzo[g,h,i]perylene | 8.7 | | 8.0 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-1 | Lab Sample ID: 240-57930-1

Matrix: Solid | Lab File ID: 51119027.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:30

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 29.52(g) | Date Analyzed: 11/19/2015 19:46

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 9.7 | | 8.0 | 0.81 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 23 | J B | 84 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 84 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 60 | 32 |
| 218-01-9 | Chrysene | 18 | | 8.0 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.79 | U | 8.0 | 0.79 |
| 132-64-9 | Dibenzofuran | 14 | J | 60 | 0.79 |
| 84-66-2 | Diethyl phthalate | 19 | U | 84 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 84 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 84 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.4 | U | 84 | 9.4 |
| 206-44-0 | Fluoranthene | 27 | | 8.0 | 0.66 |
| 86-73-7 | Fluorene | 4.8 | J | 8.0 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 8.0 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.7 | U | 60 | 6.7 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.7 | U | 390 | 9.7 |
| 67-72-1 | Hexachloroethane | 11 | U | 60 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 7.3 | J | 8.0 | 0.42 |
| 78-59-1 | Isophorone | 16 | U | 60 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.5 | U | 60 | 7.5 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 60 | 25 |
| 91-20-3 | Naphthalene | 7.4 | J | 8.0 | 0.98 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 20 | | 8.0 | 0.87 |
| 108-95-2 | Phenol | 8.7 | U | 60 | 8.7 |
| 129-00-0 | Pyrene | 23 | | 8.0 | 0.53 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-1 | Lab Sample ID: 240-57930-1

Matrix: Solid | Lab File ID: 51119027.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:30

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 29.52(g) | Date Analyzed: 11/19/2015 19:46

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 44 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 44 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 42 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 43 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 44 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 39 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: 51119026.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:35

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 29.77(g) | Date Analyzed: 11/19/2015 19:23

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 7.2 | J | 60 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.54 | U | 60 | 0.54 |
| 95-57-8 | 2-Chlorophenol | 9.8 | U | 60 | 9.8 |
| 91-57-6 | 2-Methylnaphthalene | 85 | | 7.9 | 0.60 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.9 | U | 60 | 9.9 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 60 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 60 | 15 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 10 | | 7.9 | 0.91 |
| 208-96-8 | Acenaphthylene | 4.6 | J | 7.9 | 0.42 |
| 98-86-2 | Acetophenone | 24 | J | 120 | 11 |
| 120-12-7 | Anthracene | 7.4 | J | 7.9 | 0.93 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 23 | J | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 14 | | 7.9 | 0.75 |
| 50-32-8 | Benzo[a]pyrene | 12 | | 7.9 | 0.76 |
| 205-99-2 | Benzo[b]fluoranthene | 17 | | 7.9 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 10 | | 7.9 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: 51119026.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:35

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 29.77(g) | Date Analyzed: 11/19/2015 19:23

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 8.9 | | 7.9 | 0.81 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 31 | J B | 83 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 83 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 60 | 32 |
| 218-01-9 | Chrysene | 20 | | 7.9 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.79 | U | 7.9 | 0.79 |
| 132-64-9 | Dibenzofuran | 11 | J | 60 | 0.79 |
| 84-66-2 | Diethyl phthalate | 19 | U | 83 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 83 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 83 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.4 | U | 83 | 9.4 |
| 206-44-0 | Fluoranthene | 30 | | 7.9 | 0.66 |
| 86-73-7 | Fluorene | 14 | | 7.9 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.9 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.7 | U | 60 | 6.7 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.7 | U | 390 | 9.7 |
| 67-72-1 | Hexachloroethane | 11 | U | 60 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 9.0 | | 7.9 | 0.42 |
| 78-59-1 | Isophorone | 15 | U | 60 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.5 | U | 60 | 7.5 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 60 | 25 |
| 91-20-3 | Naphthalene | 51 | | 7.9 | 0.98 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 35 | | 7.9 | 0.87 |
| 108-95-2 | Phenol | 8.7 | U | 60 | 8.7 |
| 129-00-0 | Pyrene | 28 | | 7.9 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: 51119026.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:35

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 29.77(g) | Date Analyzed: 11/19/2015 19:23

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 76 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 72 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 68 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 72 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 72 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 58 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-3 | Lab Sample ID: 240-57930-3

Matrix: Solid | Lab File ID: 51119030.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:40

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.26(g) | Date Analyzed: 11/19/2015 20:57

Con. Extract Vol.: 2(mL) | Dilution Factor: 2.5

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 85 | J | 150 | 10 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 28 | U | 290 | 28 |
| 95-95-4 | 2,4,5-Trichlorophenol | 73 | U | 440 | 73 |
| 88-06-2 | 2,4,6-Trichlorophenol | 26 | U | 440 | 26 |
| 120-83-2 | 2,4-Dichlorophenol | 59 | U | 440 | 59 |
| 105-67-9 | 2,4-Dimethylphenol | 59 | U | 440 | 59 |
| 51-28-5 | 2,4-Dinitrophenol | 62 | U | 970 | 62 |
| 121-14-2 | 2,4-Dinitrotoluene | 50 | U | 590 | 50 |
| 606-20-2 | 2,6-Dinitrotoluene | 62 | U | 590 | 62 |
| 91-58-7 | 2-Chloronaphthalene | 1.3 | U | 150 | 1.3 |
| 95-57-8 | 2-Chlorophenol | 24 | U | 150 | 24 |
| 91-57-6 | 2-Methylnaphthalene | 360 | | 20 | 1.5 |
| 95-48-7 | 2-Methylphenol | 32 | U | 590 | 32 |
| 88-74-4 | 2-Nitroaniline | 27 | U | 590 | 27 |
| 88-75-5 | 2-Nitrophenol | 24 | U | 150 | 24 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 53 | U | 290 | 53 |
| 99-09-2 | 3-Nitroaniline | 47 | U | 590 | 47 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 27 | U | 440 | 27 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 38 | U | 150 | 38 |
| 59-50-7 | 4-Chloro-3-methylphenol | 62 | U | 440 | 62 |
| 106-47-8 | 4-Chloroaniline | 50 | U | 440 | 50 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 38 | U | 150 | 38 |
| 100-01-6 | 4-Nitroaniline | 76 | U | 590 | 76 |
| 100-02-7 | 4-Nitrophenol | 50 | U | 970 | 50 |
| 83-32-9 | Acenaphthene | 350 | | 20 | 2.2 |
| 208-96-8 | Acenaphthylene | 28 | | 20 | 1.0 |
| 98-86-2 | Acetophenone | 27 | U | 290 | 27 |
| 120-12-7 | Anthracene | 590 | | 20 | 2.3 |
| 1912-24-9 | Atrazine | 27 | U | 590 | 27 |
| 100-52-7 | Benzaldehyde | 35 | U | 290 | 35 |
| 56-55-3 | Benzo[a]anthracene | 960 | | 20 | 1.8 |
| 50-32-8 | Benzo[a]pyrene | 700 | | 20 | 1.9 |
| 205-99-2 | Benzo[b]fluoranthene | 1100 | | 20 | 1.7 |
| 191-24-2 | Benzo[g,h,i]perylene | 240 | | 20 | 1.0 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-3 | Lab Sample ID: 240-57930-3

Matrix: Solid | Lab File ID: 51119030.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:40

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.26(g) | Date Analyzed: 11/19/2015 20:57

Con. Extract Vol.: 2(mL) | Dilution Factor: 2.5

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 440 | | 20 | 2.0 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 64 | U | 290 | 64 |
| 111-44-4 | Bis(2-chloroethyl)ether | 5.9 | U | 290 | 5.9 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 56 | U | 210 | 56 |
| 85-68-7 | Butyl benzyl phthalate | 29 | U | 210 | 29 |
| 105-60-2 | Caprolactam | 110 | U | 970 | 110 |
| 86-74-8 | Carbazole | 150 | | 150 | 79 |
| 218-01-9 | Chrysene | 1000 | | 20 | 3.2 |
| 53-70-3 | Dibenz(a,h)anthracene | 69 | | 20 | 1.9 |
| 132-64-9 | Dibenzofuran | 330 | | 150 | 1.9 |
| 84-66-2 | Diethyl phthalate | 47 | U | 210 | 47 |
| 131-11-3 | Dimethyl phthalate | 50 | U | 210 | 50 |
| 84-74-2 | Di-n-butyl phthalate | 44 | U | 210 | 44 |
| 117-84-0 | Di-n-octyl phthalate | 23 | U | 210 | 23 |
| 206-44-0 | Fluoranthene | 2500 | | 20 | 1.6 |
| 86-73-7 | Fluorene | 350 | | 20 | 1.6 |
| 118-74-1 | Hexachlorobenzene | 6.2 | U | 20 | 6.2 |
| 87-68-3 | Hexachlorobutadiene | 16 | U | 150 | 16 |
| 77-47-4 | Hexachlorocyclopentadiene | 24 | U | 970 | 24 |
| 67-72-1 | Hexachloroethane | 26 | U | 150 | 26 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 250 | | 20 | 1.0 |
| 78-59-1 | Isophorone | 38 | U | 150 | 38 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 18 | U | 150 | 18 |
| 86-30-6 | N-Nitrosodiphenylamine | 62 | U | 150 | 62 |
| 91-20-3 | Naphthalene | 460 | | 20 | 2.4 |
| 98-95-3 | Nitrobenzene | 6.4 | U | 290 | 6.4 |
| 87-86-5 | Pentachlorophenol | 27 | U | 440 | 27 |
| 85-01-8 | Phenanthrene | 1900 | | 20 | 2.1 |
| 108-95-2 | Phenol | 21 | U | 150 | 21 |
| 129-00-0 | Pyrene | 1800 | | 20 | 1.3 |
| 15831-10-4 | 3 & 4 Methylphenol | 59 | U | 1200 | 59 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-3 | Lab Sample ID: 240-57930-3

Matrix: Solid | Lab File ID: 51119030.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:40

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.26(g) | Date Analyzed: 11/19/2015 20:57

Con. Extract Vol.: 2(mL) | Dilution Factor: 2.5

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 76 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 82 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 71 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 80 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 77 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 65 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-4 | Lab Sample ID: 240-57930-4

Matrix: Solid | Lab File ID: 51119029.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:45

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.32(g) | Date Analyzed: 11/19/2015 20:34

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 48 | J | 60 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.54 | U | 60 | 0.54 |
| 95-57-8 | 2-Chlorophenol | 9.8 | U | 60 | 9.8 |
| 91-57-6 | 2-Methylnaphthalene | 220 | | 7.9 | 0.60 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.9 | U | 60 | 9.9 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 60 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 60 | 15 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 110 | | 7.9 | 0.90 |
| 208-96-8 | Acenaphthylene | 23 | | 7.9 | 0.42 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 160 | | 7.9 | 0.93 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 17 | J | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 200 | | 7.9 | 0.75 |
| 50-32-8 | Benzo[a]pyrene | 100 | | 7.9 | 0.76 |
| 205-99-2 | Benzo[b]fluoranthene | 240 | | 7.9 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 41 | | 7.9 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-4 | Lab Sample ID: 240-57930-4

Matrix: Solid | Lab File ID: 51119029.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:45

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.32(g) | Date Analyzed: 11/19/2015 20:34

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 71 | | 7.9 | 0.81 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 33 | J B | 83 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 83 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 60 | 32 |
| 218-01-9 | Chrysene | 260 | | 7.9 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 19 | | 7.9 | 0.79 |
| 132-64-9 | Dibenzofuran | 170 | | 60 | 0.79 |
| 84-66-2 | Diethyl phthalate | 19 | U | 83 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 83 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 83 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.4 | U | 83 | 9.4 |
| 206-44-0 | Fluoranthene | 560 | | 7.9 | 0.65 |
| 86-73-7 | Fluorene | 120 | | 7.9 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.9 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.7 | U | 60 | 6.7 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.6 | U | 390 | 9.6 |
| 67-72-1 | Hexachloroethane | 11 | U | 60 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 40 | | 7.9 | 0.42 |
| 78-59-1 | Isophorone | 15 | U | 60 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.5 | U | 60 | 7.5 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 60 | 25 |
| 91-20-3 | Naphthalene | 290 | | 7.9 | 0.98 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 550 | | 7.9 | 0.87 |
| 108-95-2 | Phenol | 8.7 | U | 60 | 8.7 |
| 129-00-0 | Pyrene | 420 | | 7.9 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-4 | Lab Sample ID: 240-57930-4

Matrix: Solid | Lab File ID: 51119029.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:45

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.32(g) | Date Analyzed: 11/19/2015 20:34

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 69 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 67 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 62 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 66 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 66 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 56 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 | Lab Sample ID: 240-57930-5

Matrix: Solid | Lab File ID: 51119014.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.24(g) | Date Analyzed: 11/19/2015 14:37

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 59 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 29 | U | 180 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10 | U | 180 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 23 | U | 180 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 23 | U | 180 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 230 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.53 | U | 59 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 9.6 | U | 59 | 9.6 |
| 91-57-6 | 2-Methylnaphthalene | 0.59 | U | 7.8 | 0.59 |
| 95-48-7 | 2-Methylphenol | 13 | U | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.7 | U | 59 | 9.7 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 59 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 59 | 15 |
| 100-01-6 | 4-Nitroaniline | 30 | U | 230 | 30 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.89 | U | 7.8 | 0.89 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.8 | 0.41 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.91 | U | 7.8 | 0.91 |
| 1912-24-9 | Atrazine | 11 | U | 230 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.74 | U | 7.8 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 0.75 | U | 7.8 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 0.69 | U | 7.8 | 0.69 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.41 | U | 7.8 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 | Lab Sample ID: 240-57930-5

Matrix: Solid | Lab File ID: 51119014.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.24(g) | Date Analyzed: 11/19/2015 14:37

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.80 | U | 7.8 | 0.80 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.3 | U | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 25 | J B | 82 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 82 | 12 |
| 105-60-2 | Caprolactam | 43 | U | 390 | 43 |
| 86-74-8 | Carbazole | 32 | U | 59 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 7.8 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.77 | U | 7.8 | 0.77 |
| 132-64-9 | Dibenzofuran | 0.77 | U | 59 | 0.77 |
| 84-66-2 | Diethyl phthalate | 19 | U | 82 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 82 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 82 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.3 | U | 82 | 9.3 |
| 206-44-0 | Fluoranthene | 0.64 | U | 7.8 | 0.64 |
| 86-73-7 | Fluorene | 0.62 | U | 7.8 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.8 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.6 | U | 59 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.5 | U | 390 | 9.5 |
| 67-72-1 | Hexachloroethane | 11 | U | 59 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.41 | U | 7.8 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 59 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.4 | U | 59 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 59 | 25 |
| 91-20-3 | Naphthalene | 0.96 | U | 7.8 | 0.96 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.86 | U | 7.8 | 0.86 |
| 108-95-2 | Phenol | 8.6 | U | 59 | 8.6 |
| 129-00-0 | Pyrene | 0.52 | U | 7.8 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 23 | U | 470 | 23 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 | Lab Sample ID: 240-57930-5

Matrix: Solid | Lab File ID: 51119014.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.24(g) | Date Analyzed: 11/19/2015 14:37

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 86 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 80 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 70 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 78 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 79 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 56 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-2 | Lab Sample ID: 240-57930-6

Matrix: Solid | Lab File ID: 51119017.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:05

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.05(g) | Date Analyzed: 11/19/2015 15:48

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 59 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.53 | U | 59 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 9.7 | U | 59 | 9.7 |
| 91-57-6 | 2-Methylnaphthalene | 0.59 | U | 7.9 | 0.59 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.8 | U | 59 | 9.8 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 59 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 59 | 15 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.90 | U | 7.9 | 0.90 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.9 | 0.41 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.92 | U | 7.9 | 0.92 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.74 | U | 7.9 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 0.76 | U | 7.9 | 0.76 |
| 205-99-2 | Benzo[b]fluoranthene | 0.70 | U | 7.9 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.41 | U | 7.9 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-2 | Lab Sample ID: 240-57930-6

Matrix: Solid | Lab File ID: 51119017.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:05

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.05(g) | Date Analyzed: 11/19/2015 15:48

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.80 | U | 7.9 | 0.80 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 38 | J B | 83 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 83 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 59 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 7.9 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.78 | U | 7.9 | 0.78 |
| 132-64-9 | Dibenzofuran | 0.78 | U | 59 | 0.78 |
| 84-66-2 | Diethyl phthalate | 19 | U | 83 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 83 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 83 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.3 | U | 83 | 9.3 |
| 206-44-0 | Fluoranthene | 0.65 | U | 7.9 | 0.65 |
| 86-73-7 | Fluorene | 0.63 | U | 7.9 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.9 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.6 | U | 59 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.6 | U | 390 | 9.6 |
| 67-72-1 | Hexachloroethane | 11 | U | 59 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.41 | U | 7.9 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 59 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.4 | U | 59 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 59 | 25 |
| 91-20-3 | Naphthalene | 0.97 | U | 7.9 | 0.97 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.86 | U | 7.9 | 0.86 |
| 108-95-2 | Phenol | 8.6 | U | 59 | 8.6 |
| 129-00-0 | Pyrene | 0.52 | U | 7.9 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 470 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-2 | Lab Sample ID: 240-57930-6

Matrix: Solid | Lab File ID: 51119017.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:05

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.05(g) | Date Analyzed: 11/19/2015 15:48

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 68 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 68 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 64 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 68 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 67 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 55 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-3 | Lab Sample ID: 240-57930-7

Matrix: Solid | Lab File ID: 51119018.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:10

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.21(g) | Date Analyzed: 11/19/2015 16:12

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 19.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.3 | U | 61 | 4.3 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 12 | U | 120 | 12 |
| 95-95-4 | 2,4,5-Trichlorophenol | 31 | U | 180 | 31 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 25 | U | 180 | 25 |
| 105-67-9 | 2,4-Dimethylphenol | 25 | U | 180 | 25 |
| 51-28-5 | 2,4-Dinitrophenol | 26 | U | 400 | 26 |
| 121-14-2 | 2,4-Dinitrotoluene | 21 | U | 250 | 21 |
| 606-20-2 | 2,6-Dinitrotoluene | 26 | U | 250 | 26 |
| 91-58-7 | 2-Chloronaphthalene | 0.55 | U | 61 | 0.55 |
| 95-57-8 | 2-Chlorophenol | 10 | U | 61 | 10 |
| 91-57-6 | 2-Methylnaphthalene | 0.61 | U | 8.2 | 0.61 |
| 95-48-7 | 2-Methylphenol | 13 | U | 250 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 250 | 11 |
| 88-75-5 | 2-Nitrophenol | 10 | U | 61 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 20 | U | 250 | 20 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 61 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 26 | U | 180 | 26 |
| 106-47-8 | 4-Chloroaniline | 21 | U | 180 | 21 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 61 | 16 |
| 100-01-6 | 4-Nitroaniline | 32 | U | 250 | 32 |
| 100-02-7 | 4-Nitrophenol | 21 | U | 400 | 21 |
| 83-32-9 | Acenaphthene | 0.93 | U | 8.2 | 0.93 |
| 208-96-8 | Acenaphthylene | 0.43 | U | 8.2 | 0.43 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.96 | U | 8.2 | 0.96 |
| 1912-24-9 | Atrazine | 11 | U | 250 | 11 |
| 100-52-7 | Benzaldehyde | 45 | J | 120 | 15 |
| 56-55-3 | Benzo[a]anthracene | 0.77 | U | 8.2 | 0.77 |
| 50-32-8 | Benzo[a]pyrene | 0.78 | U | 8.2 | 0.78 |
| 205-99-2 | Benzo[b]fluoranthene | 0.72 | U | 8.2 | 0.72 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.43 | U | 8.2 | 0.43 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-3 | Lab Sample ID: 240-57930-7

Matrix: Solid | Lab File ID: 51119018.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:10

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.21(g) | Date Analyzed: 11/19/2015 16:12

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 19.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.83 | U | 8.2 | 0.83 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 27 | U | 120 | 27 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.5 | U | 120 | 2.5 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 55 | J B | 86 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 86 | 12 |
| 105-60-2 | Caprolactam | 45 | U | 400 | 45 |
| 86-74-8 | Carbazole | 33 | U | 61 | 33 |
| 218-01-9 | Chrysene | 1.3 | U | 8.2 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.81 | U | 8.2 | 0.81 |
| 132-64-9 | Dibenzofuran | 0.81 | U | 61 | 0.81 |
| 84-66-2 | Diethyl phthalate | 20 | U | 86 | 20 |
| 131-11-3 | Dimethyl phthalate | 21 | U | 86 | 21 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 86 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.7 | U | 86 | 9.7 |
| 206-44-0 | Fluoranthene | 0.67 | U | 8.2 | 0.67 |
| 86-73-7 | Fluorene | 0.65 | U | 8.2 | 0.65 |
| 118-74-1 | Hexachlorobenzene | 2.6 | U | 8.2 | 2.6 |
| 87-68-3 | Hexachlorobutadiene | 6.9 | U | 61 | 6.9 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.9 | U | 400 | 9.9 |
| 67-72-1 | Hexachloroethane | 11 | U | 61 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.43 | U | 8.2 | 0.43 |
| 78-59-1 | Isophorone | 16 | U | 61 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.7 | U | 61 | 7.7 |
| 86-30-6 | N-Nitrosodiphenylamine | 26 | U | 61 | 26 |
| 91-20-3 | Naphthalene | 1.0 | U | 8.2 | 1.0 |
| 98-95-3 | Nitrobenzene | 2.7 | U | 120 | 2.7 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.90 | U | 8.2 | 0.90 |
| 108-95-2 | Phenol | 9.0 | U | 61 | 9.0 |
| 129-00-0 | Pyrene | 0.54 | U | 8.2 | 0.54 |
| 15831-10-4 | 3 & 4 Methylphenol | 25 | U | 490 | 25 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-3 | Lab Sample ID: 240-57930-7

Matrix: Solid | Lab File ID: 51119018.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:10

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.21(g) | Date Analyzed: 11/19/2015 16:12

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 19.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 66 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 67 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 59 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 66 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 63 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 40 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-4 | Lab Sample ID: 240-57930-8

Matrix: Solid | Lab File ID: 51119019.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:15

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 29.80(g) | Date Analyzed: 11/19/2015 16:36

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 60 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 400 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.54 | U | 60 | 0.54 |
| 95-57-8 | 2-Chlorophenol | 9.8 | U | 60 | 9.8 |
| 91-57-6 | 2-Methylnaphthalene | 0.60 | U | 8.0 | 0.60 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 10 | U | 60 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 60 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 60 | 16 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 400 | 20 |
| 83-32-9 | Acenaphthene | 0.91 | U | 8.0 | 0.91 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 8.0 | 0.42 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.94 | U | 8.0 | 0.94 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.76 | U | 8.0 | 0.76 |
| 50-32-8 | Benzo[a]pyrene | 0.77 | U | 8.0 | 0.77 |
| 205-99-2 | Benzo[b]fluoranthene | 0.71 | U | 8.0 | 0.71 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.42 | U | 8.0 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-4 | Lab Sample ID: 240-57930-8

Matrix: Solid | Lab File ID: 51119019.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:15

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 29.80(g) | Date Analyzed: 11/19/2015 16:36

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.82 | U | 8.0 | 0.82 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 30 | J B | 84 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 84 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 400 | 44 |
| 86-74-8 | Carbazole | 32 | U | 60 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 8.0 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.79 | U | 8.0 | 0.79 |
| 132-64-9 | Dibenzofuran | 0.79 | U | 60 | 0.79 |
| 84-66-2 | Diethyl phthalate | 19 | U | 84 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 84 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 84 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.5 | U | 84 | 9.5 |
| 206-44-0 | Fluoranthene | 0.66 | U | 8.0 | 0.66 |
| 86-73-7 | Fluorene | 0.64 | U | 8.0 | 0.64 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 8.0 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.7 | U | 60 | 6.7 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.7 | U | 400 | 9.7 |
| 67-72-1 | Hexachloroethane | 11 | U | 60 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.42 | U | 8.0 | 0.42 |
| 78-59-1 | Isophorone | 16 | U | 60 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.6 | U | 60 | 7.6 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 60 | 25 |
| 91-20-3 | Naphthalene | 0.98 | U | 8.0 | 0.98 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.88 | U | 8.0 | 0.88 |
| 108-95-2 | Phenol | 8.8 | U | 60 | 8.8 |
| 129-00-0 | Pyrene | 0.53 | U | 8.0 | 0.53 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-4 | Lab Sample ID: 240-57930-8

Matrix: Solid | Lab File ID: 51119019.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:15

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 29.80(g) | Date Analyzed: 11/19/2015 16:36

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 74 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 75 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 66 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 73 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 71 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 54 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-5 | Lab Sample ID: 240-57930-9

Matrix: Solid | Lab File ID: 51119020.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:20

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 29.71(g) | Date Analyzed: 11/19/2015 17:00

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.2 | U | 61 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 12 | U | 120 | 12 |
| 95-95-4 | 2,4,5-Trichlorophenol | 30 | U | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 400 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 21 | U | 240 | 21 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.55 | U | 61 | 0.55 |
| 95-57-8 | 2-Chlorophenol | 9.9 | U | 61 | 9.9 |
| 91-57-6 | 2-Methylnaphthalene | 0.61 | U | 8.1 | 0.61 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 10 | U | 61 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 61 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 21 | U | 180 | 21 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 61 | 16 |
| 100-01-6 | 4-Nitroaniline | 32 | U | 240 | 32 |
| 100-02-7 | 4-Nitrophenol | 21 | U | 400 | 21 |
| 83-32-9 | Acenaphthene | 0.92 | U | 8.1 | 0.92 |
| 208-96-8 | Acenaphthylene | 0.42 | U | 8.1 | 0.42 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.95 | U | 8.1 | 0.95 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 15 | U | 120 | 15 |
| 56-55-3 | Benzo[a]anthracene | 0.76 | U | 8.1 | 0.76 |
| 50-32-8 | Benzo[a]pyrene | 0.78 | U | 8.1 | 0.78 |
| 205-99-2 | Benzo[b]fluoranthene | 0.72 | U | 8.1 | 0.72 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.42 | U | 8.1 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-5 | Lab Sample ID: 240-57930-9

Matrix: Solid | Lab File ID: 51119020.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:20

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 29.71(g) | Date Analyzed: 11/19/2015 17:00

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.82 | U | 8.1 | 0.82 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 27 | U | 120 | 27 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 62 | J B | 85 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 85 | 12 |
| 105-60-2 | Caprolactam | 45 | U | 400 | 45 |
| 86-74-8 | Carbazole | 33 | U | 61 | 33 |
| 218-01-9 | Chrysene | 1.3 | U | 8.1 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.80 | U | 8.1 | 0.80 |
| 132-64-9 | Dibenzofuran | 0.80 | U | 61 | 0.80 |
| 84-66-2 | Diethyl phthalate | 19 | U | 85 | 19 |
| 131-11-3 | Dimethyl phthalate | 21 | U | 85 | 21 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 85 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.6 | U | 85 | 9.6 |
| 206-44-0 | Fluoranthene | 0.67 | U | 8.1 | 0.67 |
| 86-73-7 | Fluorene | 0.64 | U | 8.1 | 0.64 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 8.1 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.8 | U | 61 | 6.8 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.8 | U | 400 | 9.8 |
| 67-72-1 | Hexachloroethane | 11 | U | 61 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.42 | U | 8.1 | 0.42 |
| 78-59-1 | Isophorone | 16 | U | 61 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.6 | U | 61 | 7.6 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 61 | 25 |
| 91-20-3 | Naphthalene | 0.99 | U | 8.1 | 0.99 |
| 98-95-3 | Nitrobenzene | 2.7 | U | 120 | 2.7 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.88 | U | 8.1 | 0.88 |
| 108-95-2 | Phenol | 8.8 | U | 61 | 8.8 |
| 129-00-0 | Pyrene | 0.53 | U | 8.1 | 0.53 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 480 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-5 | Lab Sample ID: 240-57930-9

Matrix: Solid | Lab File ID: 51119020.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:20

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 29.71(g) | Date Analyzed: 11/19/2015 17:00

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 73 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 69 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 60 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 68 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 67 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 41 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-6 | Lab Sample ID: 240-57930-10

Matrix: Solid | Lab File ID: 51119021.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:25

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.14(g) | Date Analyzed: 11/19/2015 17:23

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 59 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 29 | U | 180 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10 | U | 180 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 23 | U | 180 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 23 | U | 180 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 230 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.53 | U | 59 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 9.6 | U | 59 | 9.6 |
| 91-57-6 | 2-Methylnaphthalene | 0.59 | U | 7.8 | 0.59 |
| 95-48-7 | 2-Methylphenol | 13 | U | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.7 | U | 59 | 9.7 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 59 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 59 | 15 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 230 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.89 | U | 7.8 | 0.89 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.8 | 0.41 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.92 | U | 7.8 | 0.92 |
| 1912-24-9 | Atrazine | 11 | U | 230 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 0.74 | U | 7.8 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 0.75 | U | 7.8 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 0.69 | U | 7.8 | 0.69 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.41 | U | 7.8 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-6 | Lab Sample ID: 240-57930-10

Matrix: Solid | Lab File ID: 51119021.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:25

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.14(g) | Date Analyzed: 11/19/2015 17:23

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.80 | U | 7.8 | 0.80 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.3 | U | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 31 | J B | 82 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 82 | 12 |
| 105-60-2 | Caprolactam | 43 | U | 390 | 43 |
| 86-74-8 | Carbazole | 32 | U | 59 | 32 |
| 218-01-9 | Chrysene | 1.3 | U | 7.8 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.77 | U | 7.8 | 0.77 |
| 132-64-9 | Dibenzofuran | 0.77 | U | 59 | 0.77 |
| 84-66-2 | Diethyl phthalate | 19 | U | 82 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 82 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 82 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.3 | U | 82 | 9.3 |
| 206-44-0 | Fluoranthene | 0.65 | U | 7.8 | 0.65 |
| 86-73-7 | Fluorene | 0.62 | U | 7.8 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.8 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.6 | U | 59 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.5 | U | 390 | 9.5 |
| 67-72-1 | Hexachloroethane | 11 | U | 59 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.41 | U | 7.8 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 59 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.4 | U | 59 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 59 | 25 |
| 91-20-3 | Naphthalene | 0.96 | U | 7.8 | 0.96 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 0.86 | U | 7.8 | 0.86 |
| 108-95-2 | Phenol | 8.6 | U | 59 | 8.6 |
| 129-00-0 | Pyrene | 0.52 | U | 7.8 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 23 | U | 470 | 23 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-6 | Lab Sample ID: 240-57930-10

Matrix: Solid | Lab File ID: 51119021.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:25

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.14(g) | Date Analyzed: 11/19/2015 17:23

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 70 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 69 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 61 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 68 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 66 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 53 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP | Lab Sample ID: 240-57930-11

Matrix: Solid | Lab File ID: 51119028.D

Analysis Method: 8270D | Date Collected: 11/13/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.46(g) | Date Analyzed: 11/19/2015 20:10

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 58 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 29 | U | 180 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10 | U | 180 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 23 | U | 180 | 23 |
| 105-67-9 | 2,4-Dimethylphenol | 23 | U | 180 | 23 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 230 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 230 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.53 | U | 58 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 9.6 | U | 58 | 9.6 |
| 91-57-6 | 2-Methylnaphthalene | 7.8 | | 7.8 | 0.58 |
| 95-48-7 | 2-Methylphenol | 13 | U | 230 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 230 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.7 | U | 58 | 9.7 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 230 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 58 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 58 | 15 |
| 100-01-6 | 4-Nitroaniline | 30 | U | 230 | 30 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 0.89 | U | 7.8 | 0.89 |
| 208-96-8 | Acenaphthylene | 4.0 | J | 7.8 | 0.41 |
| 98-86-2 | Acetophenone | 17 | J | 120 | 11 |
| 120-12-7 | Anthracene | 5.4 | J | 7.8 | 0.91 |
| 1912-24-9 | Atrazine | 11 | U | 230 | 11 |
| 100-52-7 | Benzaldehyde | 25 | J | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 16 | | 7.8 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 17 | | 7.8 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 30 | | 7.8 | 0.69 |
| 191-24-2 | Benzo[g,h,i]perylene | 10 | | 7.8 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP | Lab Sample ID: 240-57930-11

Matrix: Solid | Lab File ID: 51119028.D

Analysis Method: 8270D | Date Collected: 11/13/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.46(g) | Date Analyzed: 11/19/2015 20:10

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 9.9 | | 7.8 | 0.79 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.3 | U | 120 | 2.3 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 22 | U | 82 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 82 | 12 |
| 105-60-2 | Caprolactam | 43 | U | 390 | 43 |
| 86-74-8 | Carbazole | 32 | U | 58 | 32 |
| 218-01-9 | Chrysene | 22 | | 7.8 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.77 | U | 7.8 | 0.77 |
| 132-64-9 | Dibenzofuran | 16 | J | 58 | 0.77 |
| 84-66-2 | Diethyl phthalate | 19 | U | 82 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 82 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 18 | U | 82 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.2 | U | 82 | 9.2 |
| 206-44-0 | Fluoranthene | 37 | | 7.8 | 0.64 |
| 86-73-7 | Fluorene | 6.1 | J | 7.8 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.8 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.5 | U | 58 | 6.5 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.5 | U | 390 | 9.5 |
| 67-72-1 | Hexachloroethane | 11 | U | 58 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 9.2 | | 7.8 | 0.41 |
| 78-59-1 | Isophorone | 15 | U | 58 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.4 | U | 58 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 58 | 25 |
| 91-20-3 | Naphthalene | 12 | | 7.8 | 0.96 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 27 | | 7.8 | 0.85 |
| 108-95-2 | Phenol | 8.5 | U | 58 | 8.5 |
| 129-00-0 | Pyrene | 30 | | 7.8 | 0.51 |
| 15831-10-4 | 3 & 4 Methylphenol | 23 | U | 470 | 23 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP | Lab Sample ID: 240-57930-11

Matrix: Solid | Lab File ID: 51119028.D

Analysis Method: 8270D | Date Collected: 11/13/2015 00:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30.46(g) | Date Analyzed: 11/19/2015 20:10

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 66 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 68 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 65 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 68 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 66 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 62 | | 10-110 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35 Calibration End Date: 11/13/2015 13:36 Calibration ID: 31440

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-206629/6 | 51113006.D |
| Level 2 | STD2 240-206629/5 | 51113005.D |
| Level 3 | STD3 240-206629/4 | 51113004.D |
| Level 4 | STD4 240-206629/3 | 51113003.D |
| Level 5 | STD5 240-206629/2 | 51113002.D |
| Level 6 | STD6 240-206629/10 | 51113010.D |
| Level 7 | STD7 240-206629/9 | 51113009.D |
| Level 8 | STD8 240-206629/8 | 51113008.D |
| Level 9 | STD9 240-206629/12 | 51113012.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,4-Dioxane | 0.5800 | 0.6518<br>0.5735 | 0.6050<br>0.5469 | 0.6127<br>0.5769 | 0.5743 | Ave | | 0.5901 | | | | 5.4 | | 20.0 | | | |
| N-Nitrosodimethylamine | 0.7783 | 0.8219<br>0.7580 | 0.7811<br>0.7426 | 0.7616<br>0.7504 | 0.7567 | Ave | | 0.7688 | | | | 3.3 | | 20.0 | | | |
| Pyridine | 1.4242 | 1.4152<br>1.3894 | 1.3399<br>1.3496 | 1.3560<br>1.3946 | 1.3788 | Ave | | 1.3810 | | | | 2.2 | | 20.0 | | | |
| Benzaldehyde | 1.1864 | 1.2241<br>1.1177 | 1.1674<br>1.0779 | 1.1723<br>1.0659 | 1.1261 | Ave | | 1.1422 | | | 0.0100 | 4.8 | | 20.0 | | | |
| Phenol | 1.7826 | 1.8299<br>1.7560 | 1.7294<br>1.7559 | 1.7444<br>1.7386 | 1.6973 | Ave | | 1.7543 | | | 0.8000 | 2.2 | | 20.0 | | | |
| Aniline | 2.2200 | 2.0959<br>2.1926 | 2.1413<br>2.2045 | 2.1057<br>2.1225 | 2.1306 | Ave | | 2.1516 | | | | 2.2 | | 20.0 | | | |
| Bis(2-chloroethyl)ether | 1.4126 | 1.4209<br>1.3996 | 1.3734<br>1.3925 | 1.3811<br>1.3370 | 1.3403 | Ave | | 1.3822 | | | 0.7000 | 2.2 | | 20.0 | | | |
| 2-Chlorophenol | 1.5265 | 1.5902<br>1.5088 | 1.4322<br>1.5088 | 1.4392<br>1.4962 | 1.4547 | Ave | | 1.4946 | | | 0.8000 | 3.5 | | 20.0 | | | |
| n-Decane | 2.1436 | 2.2445<br>2.0504 | 2.0857<br>1.9668 | 2.1294<br>1.9880 | 2.0845 | Ave | | 2.0866 | | | | 4.3 | | 20.0 | | | |
| 1,3-Dichlorobenzene | 1.5862 | 1.5431<br>1.5577 | 1.5237<br>1.5489 | 1.5519<br>1.5476 | 1.4780 | Ave | | 1.5421 | | | | 2.0 | | 20.0 | | | |
| 1,4-Dichlorobenzene | 1.6043 | 1.5729<br>1.5913 | 1.5631<br>1.6042 | 1.5756<br>1.5877 | 1.5282 | Ave | | 1.5784 | | | | 1.6 | | 20.0 | | | |
| Benzyl alcohol | 0.9805 | 0.9353<br>0.9745 | 0.9482<br>0.9776 | 0.9603<br>0.9249 | 0.9370 | Ave | | 0.9548 | | | | 2.3 | | 20.0 | | | |
| 1,2-Dichlorobenzene | 1.5087 | 1.6362<br>1.5201 | 1.5091<br>1.4922 | 1.4979<br>1.4736 | 1.4511 | Ave | | 1.5111 | | | | 3.6 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35 Calibration End Date: 11/13/2015 13:36 Calibration ID: 31440

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2-Methylphenol | 1.3736 | 1.2808<br>1.3399 | 1.3577<br>1.3381 | 1.3330<br>1.2780 | 1.3186 | Ave | | 1.3275 | | | 0.7000 | 2.6 | | 20.0 | | | |
| bis (2-chloroisopropyl) ether | 2.5417 | 2.8607<br>2.3346 | 2.8373<br>2.1272 | 2.7307<br>2.1050 | 2.6422 | Ave | | 2.5224 | | | | 12.0 | | 20.0 | | | |
| Indene | 2.7102 | 2.6474<br>2.6984 | 2.5476<br>2.6781 | 2.5485<br>2.6219 | 2.5238 | Ave | | 2.6220 | | | | 2.8 | | 20.0 | | | |
| 3 & 4 Methylphenol | 1.4087 | 1.4373<br>1.4110 | 1.3951<br>1.4275 | 1.3518<br>1.3523 | 1.3358 | Ave | | 1.3899 | | | | 2.8 | | 20.0 | | | |
| N-Nitrosodi-n-propylamine | 0.9281 | 0.9711<br>0.9045 | 0.9778<br>0.9021 | 0.9138<br>0.8359 | 0.8942 | Ave | | 0.9159 | | | 0.5000 | 4.9 | | 20.0 | | | |
| Acetophenone | 2.0659 | 2.1044<br>2.0616 | 2.0352<br>2.0341 | 2.0250<br>1.9176 | 2.0397 | Ave | | 2.0354 | | | 0.0100 | 2.7 | | 20.0 | | | |
| Hexachloroethane | 0.6972 | 0.7208<br>0.6731 | 0.6963<br>0.6734 | 0.6884<br>0.6590 | 0.6970 | Ave | | 0.6881 | | | 0.3000 | 2.8 | | 20.0 | | | |
| Nitrobenzene | 0.3778 | 0.3749<br>0.3686 | 0.3707<br>0.3613 | 0.3816<br>0.3697 | 0.3774 | Ave | | 0.3728 | | | 0.2000 | 1.7 | | 20.0 | | | |
| Isophorone | 0.6799 | 0.6559<br>0.6732 | 0.6537<br>0.6739 | 0.6869<br>0.6736 | 0.6749 | Ave | | 0.6715 | | | 0.4000 | 1.7 | | 20.0 | | | |
| 2-Nitrophenol | 0.1876 | 0.1734<br>0.1859 | 0.1703<br>0.1891 | 0.1764<br>0.1773 | 0.1805 | Ave | | 0.1801 | | | 0.1000 | 3.8 | | 20.0 | | | |
| 2,4-Dimethylphenol | 0.3717 | 0.3742<br>0.3694 | 0.3797<br>0.3621 | 0.3792<br>0.3645 | 0.3749 | Ave | | 0.3720 | | | 0.2000 | 1.7 | | 20.0 | | | |
| Benzoic acid | 0.1888 | 0.1206<br>0.2118 | 0.1367<br>0.1951 | 0.1596<br>0.1911 | 0.1891 | Ave | | 0.1741 | | | | 18.2 | | 20.0 | | | |
| Bis(2-chloroethoxy)methane | 0.4049 | 0.3828<br>0.4012 | 0.4097<br>0.4105 | 0.4013<br>0.4036 | 0.4010 | Ave | | 0.4019 | | | 0.3000 | 2.1 | | 20.0 | | | |
| 2,4-Dichlorophenol | 0.2996 | 0.2683<br>0.2949 | 0.2853<br>0.3042 | 0.2956<br>0.3034 | 0.2910 | Ave | | 0.2928 | | | 0.2000 | 4.0 | | 20.0 | | | |
| 1,2,4-Trichlorobenzene | 0.3089 | 0.2973<br>0.3035 | 0.2865<br>0.3067 | 0.3112<br>0.3146 | 0.3047 | Ave | | 0.3042 | | | | 2.9 | | 20.0 | | | |
| Naphthalene | 1.1367<br>1.0464 | 0.9898<br>1.0495 | 0.9887<br>1.0832 | 1.0343<br>1.0809 | 1.0232 | Ave | | 1.0481 | | | 0.7000 | 4.5 | | 20.0 | | | |
| 4-Chloroaniline | 0.4403 | 0.3903<br>0.4489 | 0.4355<br>0.4593 | 0.4451<br>0.4446 | 0.4305 | Ave | | 0.4368 | | | 0.0100 | 4.7 | | 20.0 | | | |
| 2,6-Dichlorophenol | 0.2956 | 0.2749<br>0.2924 | 0.2831<br>0.3000 | 0.2996<br>0.2983 | 0.2935 | Ave | | 0.2922 | | | | 3.0 | | 20.0 | | | |
| Hexachlorobutadiene | 0.2016 | 0.2033<br>0.1960 | 0.2043<br>0.1966 | 0.1990<br>0.1949 | 0.2026 | Ave | | 0.1998 | | | 0.0100 | 1.8 | | 20.0 | | | |
| Caprolactam | 0.1048 | 0.1018<br>0.1037 | 0.0996<br>0.1101 | 0.0991<br>0.1069 | 0.1022 | Ave | | 0.1035 | | | 0.0100 | 3.6 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35 Calibration End Date: 11/13/2015 13:36 Calibration ID: 31440

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 4-Chloro-3-methylphenol | 0.3150 | 0.3032<br>0.3112 | 0.2950<br>0.3075 | 0.3066<br>0.3057 | 0.3153 | Ave | | 0.3075 | | | 0.2000 | 2.2 | | 20.0 | | | |
| 2-Methylnaphthalene | 0.7584<br>0.7045 | 0.6703<br>0.7087 | 0.6741<br>0.7278 | 0.6888<br>0.7205 | 0.6840 | Ave | | 0.7041 | | | 0.4000 | 4.0 | | 20.0 | | | |
| 1-Methylnaphthalene | 0.6888<br>0.6050 | 0.5890<br>0.6051 | 0.5702<br>0.6301 | 0.5932<br>0.6158 | 0.5897 | Ave | | 0.6096 | | | | 5.6 | | 20.0 | | | |
| Hexachlorocyclopentadiene | 0.3049 | 0.2789<br>0.3260 | 0.2926<br>0.3369 | 0.3285<br>0.2965 | 0.3588 | Ave | | 0.3154 | | | 0.0500 | 8.4 | | 20.0 | | | |
| 1,2,4,5-Tetrachlorobenzene | 0.5108 | 0.4766<br>0.5052 | 0.4944<br>0.5094 | 0.5125<br>0.5270 | 0.5127 | Ave | | 0.5061 | | | 0.0100 | 3.0 | | 20.0 | | | |
| 2,4,6-Trichlorophenol | 0.3656 | 0.3531<br>0.3709 | 0.3538<br>0.3706 | 0.3642<br>0.3760 | 0.3612 | Ave | | 0.3644 | | | 0.2000 | 2.2 | | 20.0 | | | |
| 2,4,5-Trichlorophenol | 0.3846 | 0.3304<br>0.3781 | 0.3662<br>0.3933 | 0.3740<br>0.3913 | 0.3740 | Ave | | 0.3740 | | | 0.2000 | 5.3 | | 20.0 | | | |
| 1,1'-Biphenyl | 1.4883<br>1.3543 | 1.3229<br>1.3801 | 1.3167<br>1.3861 | 1.3682<br>1.4058 | 1.3283 | Ave | | 1.3723 | | | 0.0100 | 3.9 | | 20.0 | | | |
| 2-Chloronaphthalene | 1.2122<br>1.1168 | 1.0507<br>1.1305 | 1.0503<br>1.1262 | 1.1313<br>1.1443 | 1.1240 | Ave | | 1.1207 | | | 0.8000 | 4.4 | | 20.0 | | | |
| 2-Nitroaniline | 0.3864 | 0.3519<br>0.3814 | 0.3758<br>0.3776 | 0.3868<br>0.3758 | 0.3903 | Ave | | 0.3783 | | | 0.0100 | 3.2 | | 20.0 | | | |
| Dimethyl phthalate | 1.2811 | 1.2965<br>1.2891 | 1.2452<br>1.3299 | 1.2854<br>1.2973 | 1.2875 | Ave | | 1.2890 | | | 0.0100 | 1.8 | | 20.0 | | | |
| 1,3-Dinitrobenzene | 0.2034 | 0.1538<br>0.2064 | 0.1745<br>0.2024 | 0.1812<br>0.2032 | 0.1940 | Ave | | 0.1899 | | | | 9.8 | | 20.0 | | | |
| 2,6-Dinitrotoluene | 0.2871 | 0.2590<br>0.2913 | 0.2843<br>0.2911 | 0.2753<br>0.2855 | 0.2906 | Ave | | 0.2830 | | | | 3.9 | | 20.0 | | | |
| Acenaphthylene | 1.6658<br>1.6993 | 1.5320<br>1.7126 | 1.6281<br>1.7529 | 1.6379<br>1.7469 | 1.6293 | Ave | | 1.6672 | | | 0.9000 | 4.2 | | 20.0 | | | |
| 3-Nitroaniline | 0.2992 | 0.2961<br>0.3084 | 0.2955<br>0.3120 | 0.3038<br>0.3002 | 0.3103 | Ave | | 0.3032 | | | 0.0100 | 2.1 | | 20.0 | | | |
| 2,4-Dinitrophenol | 0.1105 | +++++<br>0.1381 | 0.0794<br>0.1509 | 0.1056<br>+++++ | 0.1296 | Qua | 0.0601 | 0.0858 | 0.0015927 | | 0.0100 | | | | 0.9960 | | 0.9900 |
| Acenaphthene | 1.1725<br>1.0525 | 0.9890<br>1.0672 | 1.0580<br>1.0663 | 1.0432<br>1.0818 | 1.0251 | Ave | | 1.0617 | | | 0.9000 | 4.7 | | 20.0 | | | |
| 4-Nitrophenol | 0.2085 | 0.1942<br>0.2080 | 0.2128<br>0.2107 | 0.2115<br>0.2014 | 0.2261 | Ave | | 0.2092 | | | | 4.4 | | 20.0 | | | |
| 2,4-Dinitrotoluene | 0.3861 | 0.3383<br>0.3813 | 0.3314<br>0.3905 | 0.3661<br>0.3763 | 0.3809 | Ave | | 0.3688 | | | 0.2000 | 6.0 | | 20.0 | | | |
| Dibenzofuran | 1.6988<br>1.5987 | 1.5587<br>1.5673 | 1.5149<br>1.6099 | 1.5330<br>1.6047 | 1.5220 | Ave | | 1.5787 | | | 0.8000 | 3.6 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35  Calibration End Date: 11/13/2015 13:36  Calibration ID: 31440

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2,3,4,6-Tetrachlorophenol | 0.3308 | 0.3098<br>0.3326 | 0.3008<br>0.3407 | 0.3247<br>0.3326 | 0.3270 | Ave | | 0.3249 | | | 0.0100 | 4.1 | | 20.0 | | | |
| Diethyl phthalate | 1.2701 | 1.2539<br>1.2773 | 1.2952<br>1.2386 | 1.2884<br>1.2549 | 1.3024 | Ave | | 1.2726 | | | 0.0100 | 1.8 | | 20.0 | | | |
| Hexadecane | 0.6693 | 0.7033<br>0.6108 | 0.7106<br>0.5557 | 0.7074<br>0.6137 | 0.6766 | Ave | | 0.6559 | | | | 8.6 | | 20.0 | | | |
| 4-Chlorophenyl phenyl ether | 0.5913 | 0.5794<br>0.6030 | 0.6048<br>0.6145 | 0.5955<br>0.6023 | 0.5858 | Ave | | 0.5971 | | | 0.4000 | 1.9 | | 20.0 | | | |
| 4-Nitroaniline | 0.3028 | 0.2764<br>0.3063 | 0.2974<br>0.3178 | 0.3134<br>0.2963 | 0.3121 | Ave | | 0.3028 | | | 0.0100 | 4.3 | | 20.0 | | | |
| Fluorene | 1.5561<br>1.2847 | 1.1922<br>1.3236 | 1.2493<br>1.3456 | 1.2672<br>1.3403 | 1.2612 | Ave | | 1.3133 | | | 0.9000 | 7.9 | | 20.0 | | | |
| 4,6-Dinitro-2-methylphenol | 0.1083 | +++++<br>0.1251 | 0.0896<br>0.1353 | 0.1037<br>0.0964 | 0.1164 | Ave | | 0.1107 | | | 0.0100 | 14.5 | | 20.0 | | | |
| N-Nitrosodiphenylamine | 0.6062 | 0.5806<br>0.6089 | 0.5828<br>0.6313 | 0.5617<br>0.6381 | 0.5872 | Ave | | 0.5996 | | | 0.0100 | 4.4 | | 20.0 | | | |
| 1,2-Diphenylhydrazine(as Azobenzene) | 0.8281 | 0.7939<br>0.8037 | 0.8381<br>0.8095 | 0.8533<br>0.8231 | 0.8361 | Ave | | 0.8237 | | | | 2.2 | | 20.0 | | | |
| Azobenzene | 0.8288 | 0.7998<br>0.8036 | 0.8349<br>0.8095 | 0.8533<br>0.8233 | 0.8361 | Ave | | 0.8237 | | | | 2.2 | | 20.0 | | | |
| 4-Bromophenyl phenyl ether | 0.2316 | 0.2140<br>0.2319 | 0.2250<br>0.2403 | 0.2223<br>0.2370 | 0.2302 | Ave | | 0.2290 | | | 0.1000 | 3.7 | | 20.0 | | | |
| Hexachlorobenzene | 0.2871<br>0.2538 | 0.2494<br>0.2480 | 0.2530<br>0.2587 | 0.2390<br>0.2597 | 0.2473 | Ave | | 0.2551 | | | 0.1000 | 5.3 | | 20.0 | | | |
| Atrazine | 0.2333 | 0.2481<br>0.2301 | 0.2390<br>0.2292 | 0.2323<br>0.2296 | 0.2393 | Ave | | 0.2351 | | | 0.0100 | 2.8 | | 20.0 | | | |
| Pentachlorophenol | 0.1768 | 0.1442<br>0.1784 | 0.1481<br>0.1822 | 0.1633<br>0.1803 | 0.1693 | Ave | | 0.1678 | | | 0.0500 | 8.8 | | 20.0 | | | |
| n-Octadecane | 1.0559 | 1.0980<br>0.9783 | 1.1191<br>0.8727 | 1.1216<br>0.9263 | 1.1319 | Ave | | 1.0380 | | | | 9.6 | | 20.0 | | | |
| Phenanthrene | 1.2263<br>1.1418 | 1.1134<br>1.1415 | 1.0995<br>1.1991 | 1.0943<br>1.1964 | 1.1118 | Ave | | 1.1471 | | | 0.7000 | 4.2 | | 20.0 | | | |
| Anthracene | 1.2243<br>1.1516 | 1.0926<br>1.1652 | 1.1414<br>1.2289 | 1.1277<br>1.2029 | 1.1101 | Ave | | 1.1605 | | | 0.7000 | 4.2 | | 20.0 | | | |
| Carbazole | 1.0457 | 1.0752<br>1.0377 | 1.0810<br>1.1224 | 1.0543<br>1.0774 | 1.0361 | Ave | | 1.0662 | | | 0.0100 | 2.7 | | 20.0 | | | |
| Di-n-butyl phthalate | 1.4561 | 1.3958<br>1.4564 | 1.4296<br>1.4707 | 1.4048<br>1.4520 | 1.4467 | Ave | | 1.4390 | | | 0.0100 | 1.8 | | 20.0 | | | |
| Fluoranthene | 1.3132<br>1.2607 | 1.1716<br>1.2659 | 1.2250<br>1.3248 | 1.1658<br>1.2984 | 1.2532 | Ave | | 1.2532 | | | 0.6000 | 4.6 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35  Calibration End Date: 11/13/2015 13:36  Calibration ID: 31440

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Benzidine | 0.5363 | 0.3956<br>0.5795 | 0.4461<br>0.5976 | 0.4937<br>0.4961 | 0.5614 | Ave | | 0.5133 | | | | 13.5 | | 20.0 | | | |
| Pyrene | 1.2825<br>1.1398 | 1.0882<br>1.1039 | 1.0551<br>1.1376 | 1.0938<br>1.1311 | 1.0732 | Ave | | 1.1228 | | | 0.6000 | 5.9 | | 20.0 | | | |
| Butyl benzyl phthalate | 0.5639 | 0.5476<br>0.5408 | 0.5274<br>0.5413 | 0.5546<br>0.5329 | 0.5434 | Ave | | 0.5440 | | | 0.0100 | 2.1 | | 20.0 | | | |
| 3,3'-Dichlorobenzidine | 0.4080 | 0.3631<br>0.4063 | 0.3772<br>0.4230 | 0.3889<br>0.4081 | 0.3956 | Ave | | 0.3962 | | | 0.0100 | 4.9 | | 20.0 | | | |
| Bis(2-ethylhexyl) phthalate | 0.8077 | 0.7754<br>0.7684 | 0.7957<br>0.7427 | 0.7881<br>0.7630 | 0.7902 | Ave | | 0.7789 | | | 0.0100 | 2.7 | | 20.0 | | | |
| Benzo[a]anthracene | 1.2644<br>1.0828 | 1.0291<br>1.0505 | 1.0476<br>1.0923 | 1.0611<br>1.0677 | 1.0294 | Ave | | 1.0806 | | | 0.8000 | 6.7 | | 20.0 | | | |
| Chrysene | 1.1326<br>1.0034 | 0.9676<br>0.9763 | 0.9570<br>1.0083 | 0.9767<br>0.9824 | 0.9510 | Ave | | 0.9950 | | | 0.7000 | 5.5 | | 20.0 | | | |
| Di-n-octyl phthalate | 1.6551 | 1.5266<br>1.6451 | 1.5650<br>1.6721 | 1.6257<br>1.6900 | 1.5890 | Ave | | 1.6211 | | | 0.0100 | 3.5 | | 20.0 | | | |
| Benzo[b]fluoranthene | 1.2496<br>1.2002 | 1.1171<br>1.2131 | 1.0756<br>1.2289 | 1.1259<br>1.2189 | 1.1602 | Ave | | 1.1766 | | | 0.7000 | 5.1 | | 20.0 | | | |
| Benzo[k]fluoranthene | 1.3243<br>1.2239 | 1.1343<br>1.2162 | 1.1885<br>1.2577 | 1.2414<br>1.2919 | 1.1974 | Ave | | 1.2306 | | | 0.7000 | 4.6 | | 20.0 | | | |
| Benzo[a]pyrene | 1.1558<br>1.1387 | 1.0681<br>1.1582 | 1.0554<br>1.2017 | 1.1164<br>1.1780 | 1.1000 | Ave | | 1.1302 | | | 0.7000 | 4.4 | | 20.0 | | | |
| Indeno[1,2,3-cd]pyrene | 1.3217<br>1.2589 | 1.1565<br>1.2554 | 1.1606<br>1.3417 | 1.2128<br>1.2977 | 1.2299 | Ave | | 1.2484 | | | 0.5000 | 5.3 | | 20.0 | | | |
| Dibenz(a,h)anthracene | 1.0845<br>1.0450 | 0.9071<br>1.0768 | 0.9519<br>1.1527 | 0.9946<br>1.1017 | 1.0195 | Ave | | 1.0371 | | | 0.4000 | 7.4 | | 20.0 | | | |
| Benzo[g,h,i]perylene | 1.0895<br>1.0332 | 0.9955<br>1.0569 | 0.9874<br>1.1041 | 1.0518<br>1.0589 | 1.0382 | Ave | | 1.0462 | | | 0.5000 | 3.7 | | 20.0 | | | |
| 2-Fluorophenol (Surr) | 1.4168 | 1.3862<br>1.3812 | 1.4000<br>1.3367 | 1.3188<br>1.3588 | 1.2885 | Ave | | 1.3609 | | | | 3.2 | | 20.0 | | | |
| Phenol-d5 (Surr) | 1.7114 | 1.6292<br>1.6555 | 1.6082<br>1.6630 | 1.6510<br>1.6264 | 1.6328 | Ave | | 1.6472 | | | | 1.9 | | 20.0 | | | |
| Nitrobenzene-d5 (Surr) | 0.4549<br>0.4152 | 0.3856<br>0.3887 | 0.3897<br>0.3872 | 0.4124<br>0.3887 | 0.4159 | Ave | | 0.4042 | | | | 5.7 | | 20.0 | | | |
| 2-Fluorobiphenyl (Surr) | 1.3419<br>1.2271 | 1.1704<br>1.2356 | 1.1679<br>1.2420 | 1.2420<br>1.2595 | 1.2243 | Ave | | 1.2345 | | | | 4.1 | | 20.0 | | | |
| 2,4,6-Tribromophenol (Surr) | 0.1904 | 0.1762<br>0.1938 | 0.1774<br>0.1972 | 0.1726<br>0.1931 | 0.1844 | Ave | | 0.1856 | | | | 5.0 | | 20.0 | | | |
| Terphenyl-d14 (Surr) | 0.9377<br>0.8182 | 0.7913<br>0.7941 | 0.7751<br>0.8137 | 0.7944<br>0.7943 | 0.7757 | Ave | | 0.8105 | | | | 6.2 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35 Calibration End Date: 11/13/2015 13:36 Calibration ID: 31440

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-206629/6 | 51113006.D |
| Level 2 | STD2 240-206629/5 | 51113005.D |
| Level 3 | STD3 240-206629/4 | 51113004.D |
| Level 4 | STD4 240-206629/3 | 51113003.D |
| Level 5 | STD5 240-206629/2 | 51113002.D |
| Level 6 | STD6 240-206629/10 | 51113010.D |
| Level 7 | STD7 240-206629/9 | 51113009.D |
| Level 8 | STD8 240-206629/8 | 51113008.D |
| Level 9 | STD9 240-206629/12 | 51113012.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1,4-Dioxane | DCB | Ave | 242487 | 12707<br>363801 | 22045<br>552856 | 47335<br>521483 | 102317 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| N-Nitrosodimethylamine | DCB | Ave | 325392 | 16024<br>480840 | 28462<br>750792 | 58840<br>678349 | 134807 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Pyridine | DCB | Ave | 595421 | 27590<br>881420 | 48824<br>1364385 | 104766<br>1260637 | 245629 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzaldehyde | DCB | Ave | 991964 | 47731<br>1418129 | 85080<br>2179498 | 181143<br>1927001 | 401226 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Phenol | DCB | Ave | 745241 | 35675<br>1113987 | 63018<br>1775169 | 134771<br>1571572 | 302370 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Aniline | DCB | Ave | 928081 | 40861<br>1390945 | 78027<br>2228638 | 162681<br>1918607 | 379571 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Bis(2-chloroethyl)ether | DCB | Ave | 590544 | 27703<br>887912 | 50045<br>1407822 | 106700<br>1208609 | 238771 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Chlorophenol | DCB | Ave | 638153 | 31002<br>957175 | 52189<br>1525354 | 111194<br>1352489 | 259163 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| n-Decane | DCB | Ave | 896169 | 43760<br>1300739 | 76002<br>1988420 | 164515<br>1797052 | 371359 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,3-Dichlorobenzene | DCB | Ave | 663141 | 30084<br>988203 | 55522<br>1565862 | 119894<br>1398968 | 263314 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,4-Dichlorobenzene | DCB | Ave | 670689 | 30666<br>1009476 | 56958<br>1621795 | 121730<br>1435180 | 272256 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzyl alcohol | DCB | Ave | 409889 | 18234<br>618199 | 34550<br>988372 | 74188<br>836072 | 166920 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2-Dichlorobenzene | DCB | Ave | 630733 | 31899<br>964317 | 54990<br>1508520 | 115724<br>1332087 | 258518 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Methylphenol | DCB | Ave | 574266 | 24970<br>850013 | 49472<br>1352761 | 102988<br>1155228 | 234918 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35 Calibration End Date: 11/13/2015 13:36 Calibration ID: 31440

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| bis (2-chloroisopropyl) ether | DCB | Ave | 1062591 | 55773<br>1481057 | 103389<br>2150528 | 210969<br>1902780 | 470708 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Indene | DCB | Ave | 2266057 | 103230<br>3423637 | 185664<br>5415021 | 393791<br>4740216 | 899220 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 3 & 4 Methylphenol | DCB | Ave | 588910 | 28022<br>895118 | 50835<br>1443127 | 104435<br>1222420 | 237966 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| N-Nitrosodi-n-propylamine | DCB | Ave | 388013 | 18933<br>573802 | 35631<br>911950 | 70600<br>755572 | 159295 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Acetophenone | DCB | Ave | 863670 | 41028<br>1307832 | 74160<br>2056373 | 156452<br>1733452 | 363374 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachloroethane | DCB | Ave | 291479 | 14053<br>426998 | 25373<br>680793 | 53181<br>595715 | 124168 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Nitrobenzene | NPT | Ave | 658439 | 31200<br>980027 | 55607<br>1527300 | 118322<br>1305992 | 270146 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Isophorone | NPT | Ave | 1184947 | 54579<br>1789946 | 98053<br>2848963 | 212987<br>2379097 | 483047 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Nitrophenol | NPT | Ave | 327054 | 14433<br>494204 | 25541<br>799418 | 54706<br>626179 | 129185 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dimethylphenol | NPT | Ave | 647836 | 31141<br>982084 | 56951<br>1530840 | 117586<br>1287317 | 268349 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzoic acid | NPT | Ave | 657994 | 20071<br>1126513 | 41009<br>1649366 | 98993<br>1349950 | 270718 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Bis(2-chloroethoxy)methane | NPT | Ave | 705775 | 31852<br>1066825 | 61452<br>1735387 | 124444<br>1425561 | 287039 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dichlorophenol | NPT | Ave | 522159 | 22329<br>784202 | 42790<br>1286114 | 91672<br>1071599 | 208266 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2,4-Trichlorobenzene | NPT | Ave | 538428 | 24740<br>807081 | 42975<br>1296621 | 96500<br>1111072 | 218106 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Naphthalene | NPT | Ave | 19468<br>1823811 | 82370<br>2790386 | 148289<br>4579201 | 320734<br>3817926 | 732351 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Chloroaniline | NPT | Ave | 767460 | 32477<br>1193563 | 65318<br>1941823 | 138016<br>1570389 | 308128 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,6-Dichlorophenol | NPT | Ave | 515239 | 22873<br>777432 | 42458<br>1268163 | 92896<br>1053463 | 210104 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorobutadiene | NPT | Ave | 351446 | 16918<br>521242 | 30637<br>831224 | 61702<br>688288 | 145027 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Caprolactam | NPT | Ave | 365270 | 16942<br>551640 | 29887<br>930933 | 61467<br>755257 | 146226 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 4-Chloro-3-methylphenol | NPT | Ave | 549110 | 25230<br>827429 | 44252<br>1299928 | 95087<br>1079769 | 225669 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Methylnaphthalene | NPT | Ave | 12988<br>1227817 | 55776<br>1884334 | 101112<br>3076748 | 213574<br>2544802 | 489549 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35 Calibration End Date: 11/13/2015 13:36 Calibration ID: 31440

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1-Methylnaphthalene | NPT | Ave | 11797<br>1054541 | 49011<br>1608797 | 85517<br>2663612 | 183934<br>2174950 | 422088 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorocyclopentadiene | ANT | Ave | <br>310267 | 13639<br>505734 | 25689<br>852674 | 57410<br>604183 | 147706 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2,4,5-Tetrachlorobenzene | ANT | Ave | <br>519837 | 23309<br>783706 | 43411<br>1289365 | 89558<br>1073839 | 211045 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,6-Trichlorophenol | ANT | Ave | <br>372130 | 17269<br>575315 | 31063<br>938164 | 63639<br>766180 | 148673 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,5-Trichlorophenol | ANT | Ave | <br>391450 | 16161<br>586483 | 32150<br>995421 | 65361<br>797362 | 153968 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,1'-Biphenyl | ANT | Ave | 14900<br>1378317 | 64706<br>2140847 | 115610<br>3508661 | 239083<br>2864509 | 546801 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Chloronaphthalene | ANT | Ave | 12136<br>1136589 | 51392<br>1753613 | 92221<br>2850704 | 197678<br>2331795 | 462709 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Nitroaniline | ANT | Ave | <br>393286 | 17213<br>591650 | 32999<br>955860 | 67582<br>765845 | 160685 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dimethyl phthalate | ANT | Ave | <br>1303832 | 63413<br>1999692 | 109330<br>3366191 | 224618<br>2643447 | 530004 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,3-Dinitrobenzene | ANT | Ave | <br>206981 | 7523<br>320163 | 15323<br>512282 | 31658<br>414025 | 79851 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,6-Dinitrotoluene | ANT | Ave | <br>292196 | 12666<br>451929 | 24966<br>736972 | 48109<br>581863 | 119624 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Acenaphthylene | ANT | Ave | 16677<br>1729464 | 74932<br>2656651 | 142947<br>4436952 | 286219<br>3559722 | 670718 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 3-Nitroaniline | ANT | Ave | <br>304529 | 14484<br>478463 | 25948<br>789749 | 53082<br>611705 | 127718 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dinitrophenol | ANT | Qua | <br>224981 | +++++<br>428528 | 13949<br>763934 | 36919<br>+++++ | 106677 | <br>20.0 | +++++<br>30.0 | 2.00<br>40.0 | 4.00<br>+++++ | 10.0 |
| Acenaphthene | ANT | Ave | 11739<br>1071134 | 48373<br>1655487 | 92893<br>2699034 | 182296<br>2204399 | 421989 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Nitrophenol | ANT | Ave | <br>424314 | 19000<br>645189 | 37377<br>1066898 | 73905<br>820683 | 186179 | <br>20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 2,4-Dinitrotoluene | ANT | Ave | <br>392966 | 16545<br>591541 | 29095<br>988394 | 63966<br>766687 | 156804 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dibenzofuran | ANT | Ave | 17008<br>1627061 | 76236<br>2431243 | 133013<br>4075028 | 267878<br>3269978 | 626543 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,3,4,6-Tetrachlorophenol | ANT | Ave | <br>336673 | 15151<br>515984 | 26414<br>862382 | 56745<br>677769 | 134625 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Diethyl phthalate | ANT | Ave | <br>1292620 | 61332<br>1981326 | 113722<br>3135146 | 225138<br>2557069 | 536128 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexadecane | PHN | Ave | <br>1048902 | 52981<br>1498801 | 93616<br>2186134 | 190674<br>1909356 | 433277 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35 Calibration End Date: 11/13/2015 13:36 Calibration ID: 31440

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 4-Chlorophenyl phenyl ether | ANT | Ave | 601816 | 28341<br>935407 | 53101<br>1555336 | 104064<br>1227372 | 241155 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Nitroaniline | ANT | Ave | 308154 | 13517<br>475122 | 26114<br>804383 | 54767<br>603811 | 128477 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Fluorene | ANT | Ave | 15579<br>1307520 | 58311<br>2053172 | 109688<br>3405946 | 221437<br>2731175 | 519169 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4,6-Dinitro-2-methylphenol | PHN | Ave | 339317 | +++++<br>613821 | 23619<br>1064833 | 55889<br>600116 | 149073 | 20.0 | +++++<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| N-Nitrosodiphenylamine | PHN | Ave | 1900033 | 87474<br>2988214 | 153540<br>4966871 | 302792<br>3970709 | 752035 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | PHN | Ave | 1297814 | 59805<br>1972102 | 110406<br>3184726 | 229993<br>2560884 | 535383 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Azobenzene | PHN | Ave | 1298882 | 60250<br>1971871 | 109992<br>3184726 | 229993<br>2561734 | 535405 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Bromophenyl phenyl ether | PHN | Ave | 362907 | 16124<br>569068 | 29637<br>945348 | 59913<br>737353 | 147445 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorobenzene | PHN | Ave | 4510<br>397823 | 18788<br>608443 | 33327<br>1017580 | 64405<br>807944 | 158343 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Atrazine | PHN | Ave | 731343 | 37382<br>1129111 | 62983<br>1803186 | 125218<br>1428560 | 306484 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Pentachlorophenol | PHN | Ave | 554277 | 21726<br>875723 | 39025<br>1433588 | 88033<br>1121777 | 216877 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| n-Octadecane | ANT | Ave | 1074609 | 53703<br>1517540 | 98258<br>2208985 | 195987<br>1887436 | 465957 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Phenanthrene | PHN | Ave | 19267<br>1789509 | 83873<br>2800992 | 144839<br>4717241 | 294954<br>3722367 | 711946 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Anthracene | PHN | Ave | 19235<br>1804767 | 82310<br>2859224 | 150366<br>4834609 | 303944<br>3742798 | 710886 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Carbazole | PHN | Ave | 1638804 | 81000<br>2546451 | 142411<br>4415789 | 284162<br>3352067 | 663503 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Di-n-butyl phthalate | PHN | Ave | 2282096 | 105145<br>3573885 | 188333<br>5785988 | 378636<br>4517663 | 926414 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Fluoranthene | PHN | Ave | 20632<br>1975866 | 88263<br>3106427 | 161380<br>5212018 | 314215<br>4039715 | 802530 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzidine | CRY | Ave | 1926664 | 69378<br>3349409 | 139660<br>5667898 | 301859<br>3710065 | 859695 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Pyrene | CRY | Ave | 22708<br>2047232 | 95415<br>3190127 | 165176<br>5395164 | 334352<br>4229253 | 821706 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Butyl benzyl phthalate | CRY | Ave | 1012819 | 48016<br>1562675 | 82568<br>2567274 | 169546<br>1992494 | 416025 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 3,3'-Dichlorobenzidine | CRY | Ave | 1465512 | 63673<br>2348032 | 118098<br>4011733 | 237735<br>3051543 | 605753 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 206629

SDG No.:

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/13/2015 09:35 Calibration End Date: 11/13/2015 13:36 Calibration ID: 31440

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Bis(2-ethylhexyl) phthalate | CRY | Ave | 1450736 | 67985<br>2220542 | 124565<br>3522065 | 240922<br>2852931 | 605007 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[a]anthracene | CRY | Ave | 22389<br>1944809 | 90237<br>3035848 | 163998<br>5180128 | 324359<br>3991935 | 788168 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Chrysene | CRY | Ave | 20055<br>1802245 | 84839<br>2821437 | 149813<br>4781902 | 298550<br>3673097 | 728079 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Di-n-octyl phthalate | PRY | Ave | 2531103 | 112642<br>3941562 | 210278<br>6503579 | 419689<br>5083436 | 1029544 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[b]fluoranthene | PRY | Ave | 19022<br>1835386 | 82425<br>2906433 | 144524<br>4779806 | 290651<br>3666458 | 751726 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[k]fluoranthene | PRY | Ave | 20159<br>1871666 | 83696<br>2914076 | 159694<br>4891593 | 320482<br>3886014 | 775870 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[a]pyrene | PRY | Ave | 17594<br>1741452 | 78807<br>2775036 | 141805<br>4673787 | 288195<br>3543188 | 712763 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Indeno[1,2,3-cd]pyrene | PRY | Ave | 20119<br>1925202 | 85328<br>3007984 | 155947<br>5218420 | 313092<br>3903319 | 796893 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dibenz(a,h)anthracene | PRY | Ave | 16509<br>1598177 | 66932<br>2579861 | 127895<br>4483316 | 256748<br>3313767 | 660600 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[g,h,i]perylene | PRY | Ave | 16584<br>1580038 | 73449<br>2532221 | 132670<br>4294436 | 271514<br>3185151 | 672715 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Fluorophenol (Surr) | DCB | Ave | 592303 | 27025<br>876240 | 51014<br>1351392 | 101888<br>1228319 | 229547 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Phenol-d5 (Surr) | DCB | Ave | 715480 | 31763<br>1050255 | 58602<br>1681199 | 127550<br>1470164 | 290894 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Nitrobenzene-d5 (Surr) | NPT | Ave | 7791<br>723587 | 32085<br>1033503 | 58457<br>1636760 | 127882<br>1372772 | 297702 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Fluorobiphenyl (Surr) | ANT | Ave | 13435<br>1248919 | 57244<br>1916614 | 102541<br>3143694 | 217027<br>2566398 | 503998 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,6-Tribromophenol (Surr) | ANT | Ave | 193743 | 8617<br>300602 | 15573<br>499180 | 30154<br>393558 | 75894 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Terphenyl-d14 (Surr) | CRY | Ave | 16604<br>1469620 | 69384<br>2294804 | 121331<br>3859084 | 242834<br>2970004 | 593873 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Qua = Quadratic ISTD |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID (1): CCV 240-207287/2 CCVIS Instrument ID (1): A4HP10

GC Column (1): DB-624 ID: 0.18(mm) Date Analyzed (1): 11/18/2015 08:26

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 2-Fluorophenol (Surr) | 4.85 | 100.0 |
| Benzaldehyde | 5.67 | 100.0 |
| Phenol-d5 (Surr) | 5.72 | 100.0 |
| Phenol | 5.73 | 100.0 |
| Bis(2-chloroethyl)ether | 5.81 | 100.0 |
| 2-Chlorophenol | 5.88 | 100.0 |
| 2-Methylphenol | 6.27 | 100.0 |
| bis (2-chloroisopropyl) ether | 6.30 | 100.0 |
| 3 & 4 Methylphenol | 6.41 | 100.0 |
| N-Nitrosodi-n-propylamine | 6.41 | 100.0 |
| Acetophenone | 6.42 | 100.0 |
| Hexachloroethane | 6.53 | 100.0 |
| Nitrobenzene-d5 (Surr) | 6.55 | 100.0 |
| Nitrobenzene | 6.56 | 100.0 |
| Isophorone | 6.77 | 100.0 |
| 2-Nitrophenol | 6.85 | 100.0 |
| 2,4-Dimethylphenol | 6.86 | 100.0 |
| Bis(2-chloroethoxy)methane | 6.94 | 100.0 |
| 2,4-Dichlorophenol | 7.05 | 100.0 |
| Naphthalene | 7.20 | 100.0 |
| 4-Chloroaniline | 7.23 | 100.0 |
| Hexachlorobutadiene | 7.32 | 100.0 |
| Caprolactam | 7.52 | 100.0 |
| 4-Chloro-3-methylphenol | 7.63 | 100.0 |
| 2-Methylnaphthalene | 7.80 | 100.0 |
| Hexachlorocyclopentadiene | 7.94 | 100.0 |
| 2,4,6-Trichlorophenol | 8.03 | 100.0 |
| 2,4,5-Trichlorophenol | 8.06 | 100.0 |
| 2-Fluorobiphenyl (Surr) | 8.10 | 100.0 |
| 1,1'-Biphenyl | 8.19 | 100.0 |
| 2-Chloronaphthalene | 8.22 | 100.0 |
| 2-Nitroaniline | 8.28 | 100.0 |
| Dimethyl phthalate | 8.42 | 100.0 |
| 2,6-Dinitrotoluene | 8.47 | 100.0 |
| Acenaphthylene | 8.57 | 100.0 |
| 3-Nitroaniline | 8.63 | 100.0 |
| 2,4-Dinitrophenol | 8.71 | 100.0 |
| Acenaphthene | 8.71 | 100.0 |
| 4-Nitrophenol | 8.74 | 100.0 |
| 2,4-Dinitrotoluene | 8.82 | 100.0 |
| Dibenzofuran | 8.86 | 100.0 |
| Diethyl phthalate | 9.01 | 100.0 |
| 4-Chlorophenyl phenyl ether | 9.13 | 100.0 |
| 4-Nitroaniline | 9.14 | 100.0 |
| Fluorene | 9.15 | 100.0 |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID (1): CCV 240-207287/2 CCVIS Instrument ID (1): A4HP10

GC Column (1): DB-624 ID: 0.18(mm) Date Analyzed (1): 11/18/2015 08:26

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 4,6-Dinitro-2-methylphenol | 9.17 | 100.0 |
| N-Nitrosodiphenylamine | 9.23 | 100.0 |
| 2,4,6-Tribromophenol (Surr) | 9.36 | 100.0 |
| 4-Bromophenyl phenyl ether | 9.55 | 100.0 |
| Hexachlorobenzene | 9.63 | 100.0 |
| Atrazine | 9.66 | 100.0 |
| Pentachlorophenol | 9.79 | 100.0 |
| Phenanthrene | 9.97 | 100.0 |
| Anthracene | 10.01 | 100.0 |
| Carbazole | 10.14 | 100.0 |
| Di-n-butyl phthalate | 10.38 | 100.0 |
| Fluoranthene | 10.99 | 100.0 |
| Pyrene | 11.21 | 100.0 |
| Terphenyl-d14 (Surr) | 11.32 | 100.0 |
| Butyl benzyl phthalate | 11.81 | 100.0 |
| 3,3'-Dichlorobenzidine | 12.49 | 100.0 |
| Bis(2-ethylhexyl) phthalate | 12.50 | 100.0 |
| Benzo[a]anthracene | 12.55 | 100.0 |
| Chrysene | 12.60 | 100.0 |
| Di-n-octyl phthalate | 13.42 | 100.0 |
| Benzo[b]fluoranthene | 14.10 | 26.50 |
| Benzo[k]fluoranthene | 14.14 | 100.0 |
| Benzo[a]pyrene | 14.61 | 100.0 |
| Indeno[1,2,3-cd]pyrene | 16.41 | 100.0 |
| Dibenz(a,h)anthracene | 16.43 | 100.0 |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID (1): CCV 240-207541/2 CCVIS Instrument ID (1): A4HP10

GC Column (1): DB-624 ID: 0.18(mm) Date Analyzed (1): 11/19/2015 09:50

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 2-Fluorophenol (Surr) | 4.84 | 100.0 |
| Benzaldehyde | 5.66 | 100.0 |
| Phenol-d5 (Surr) | 5.71 | 100.0 |
| Phenol | 5.72 | 100.0 |
| Bis(2-chloroethyl)ether | 5.80 | 100.0 |
| 2-Chlorophenol | 5.88 | 100.0 |
| 2-Methylphenol | 6.26 | 100.0 |
| bis (2-chloroisopropyl) ether | 6.29 | 100.0 |
| 3 & 4 Methylphenol | 6.39 | 100.0 |
| N-Nitrosodi-n-propylamine | 6.40 | 100.0 |
| Acetophenone | 6.41 | 100.0 |
| Hexachloroethane | 6.52 | 100.0 |
| Nitrobenzene-d5 (Surr) | 6.54 | 100.0 |
| Nitrobenzene | 6.56 | 100.0 |
| Isophorone | 6.76 | 100.0 |
| 2-Nitrophenol | 6.83 | 100.0 |
| 2,4-Dimethylphenol | 6.86 | 100.0 |
| Bis(2-chloroethoxy)methane | 6.93 | 100.0 |
| 2,4-Dichlorophenol | 7.05 | 100.0 |
| Naphthalene | 7.19 | 100.0 |
| 4-Chloroaniline | 7.22 | 100.0 |
| Hexachlorobutadiene | 7.30 | 100.0 |
| Caprolactam | 7.51 | 100.0 |
| 4-Chloro-3-methylphenol | 7.63 | 100.0 |
| 2-Methylnaphthalene | 7.79 | 100.0 |
| Hexachlorocyclopentadiene | 7.93 | 100.0 |
| 2,4,6-Trichlorophenol | 8.02 | 100.0 |
| 2,4,5-Trichlorophenol | 8.06 | 100.0 |
| 2-Fluorobiphenyl (Surr) | 8.09 | 100.0 |
| 1,1'-Biphenyl | 8.17 | 100.0 |
| 2-Chloronaphthalene | 8.20 | 100.0 |
| 2-Nitroaniline | 8.27 | 100.0 |
| Dimethyl phthalate | 8.41 | 100.0 |
| 2,6-Dinitrotoluene | 8.46 | 100.0 |
| Acenaphthylene | 8.56 | 100.0 |
| 3-Nitroaniline | 8.61 | 100.0 |
| 2,4-Dinitrophenol | 8.70 | 100.0 |
| Acenaphthene | 8.70 | 100.0 |
| 4-Nitrophenol | 8.74 | 100.0 |
| 2,4-Dinitrotoluene | 8.81 | 100.0 |
| Dibenzofuran | 8.85 | 100.0 |
| Diethyl phthalate | 9.00 | 100.0 |
| 4-Chlorophenyl phenyl ether | 9.12 | 100.0 |
| 4-Nitroaniline | 9.13 | 100.0 |
| Fluorene | 9.14 | 100.0 |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID (1): CCV 240-207541/2 CCVIS Instrument ID (1): A4HP10

GC Column (1): DB-624 ID: 0.18(mm) Date Analyzed (1): 11/19/2015 09:50

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 4,6-Dinitro-2-methylphenol | 9.16 | 100.0 |
| N-Nitrosodiphenylamine | 9.21 | 100.0 |
| 2,4,6-Tribromophenol (Surr) | 9.34 | 100.0 |
| 4-Bromophenyl phenyl ether | 9.54 | 100.0 |
| Hexachlorobenzene | 9.63 | 100.0 |
| Atrazine | 9.65 | 100.0 |
| Pentachlorophenol | 9.78 | 100.0 |
| Phenanthrene | 9.96 | 100.0 |
| Anthracene | 10.00 | 100.0 |
| Carbazole | 10.12 | 100.0 |
| Di-n-butyl phthalate | 10.37 | 100.0 |
| Fluoranthene | 10.98 | 100.0 |
| Pyrene | 11.20 | 100.0 |
| Terphenyl-d14 (Surr) | 11.31 | 100.0 |
| Butyl benzyl phthalate | 11.79 | 100.0 |
| 3,3'-Dichlorobenzidine | 12.47 | 100.0 |
| Bis(2-ethylhexyl) phthalate | 12.48 | 100.0 |
| Benzo[a]anthracene | 12.53 | 100.0 |
| Chrysene | 12.59 | 100.0 |
| Di-n-octyl phthalate | 13.40 | 100.0 |
| Benzo[b]fluoranthene | 14.09 | 25.20 |
| Benzo[k]fluoranthene | 14.13 | 100.0 |
| Benzo[a]pyrene | 14.59 | 100.0 |
| Indeno[1,2,3-cd]pyrene | 16.39 | 100.0 |
| Dibenz(a,h)anthracene | 16.41 | 100.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-206629/11 Calibration Date: 11/13/2015 13:12

Instrument ID: A4HP10 Calib Start Date: 11/13/2015 09:35

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/13/2015 13:36

Lab File ID: 51113011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.5901 | 0.5299 | | 8.98 | 10.0 | -10.2 | 30.0 |
| N-Nitrosodimethylamine | Ave | 0.7688 | 0.6186 | | 8.05 | 10.0 | -19.5 | 30.0 |
| Pyridine | Ave | 1.381 | 1.303 | | 9.43 | 10.0 | -5.7 | 30.0 |
| Benzaldehyde | Ave | 1.142 | 1.032 | 0.0100 | 18.1 | 20.0 | -9.7 | 30.0 |
| Phenol | Ave | 1.754 | 1.651 | 0.8000 | 9.41 | 10.0 | -5.9 | 30.0 |
| Aniline | Ave | 2.152 | 1.999 | | 9.29 | 10.0 | -7.1 | 30.0 |
| Bis(2-chloroethyl)ether | Ave | 1.382 | 1.274 | 0.7000 | 9.21 | 10.0 | -7.9 | 30.0 |
| 2-Chlorophenol | Ave | 1.495 | 1.381 | 0.8000 | 9.24 | 10.0 | -7.6 | 30.0 |
| n-Decane | Ave | 2.087 | 1.904 | | 9.13 | 10.0 | -8.7 | 30.0 |
| 1,3-Dichlorobenzene | Ave | 1.542 | 1.441 | | 9.34 | 10.0 | -6.6 | 30.0 |
| 1,4-Dichlorobenzene | Ave | 1.578 | 1.472 | | 9.33 | 10.0 | -6.7 | 30.0 |
| Benzyl alcohol | Ave | 0.9548 | 0.9034 | | 9.46 | 10.0 | -5.4 | 30.0 |
| 1,2-Dichlorobenzene | Ave | 1.511 | 1.370 | | 9.06 | 10.0 | -9.4 | 30.0 |
| 2-Methylphenol | Ave | 1.327 | 1.276 | 0.7000 | 9.61 | 10.0 | -3.9 | 30.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.522 | 2.334 | | 9.25 | 10.0 | -7.5 | 30.0 |
| Indene | Ave | 2.622 | 2.310 | | 17.6 | 20.0 | -11.9 | 30.0 |
| 3 & 4 Methylphenol | Ave | 1.390 | 1.301 | | 9.36 | 10.0 | -6.4 | 30.0 |
| N-Nitrosodi-n-propylamine | Ave | 0.9159 | 0.8494 | 0.5000 | 9.27 | 10.0 | -7.3 | 30.0 |
| Acetophenone | Ave | 2.035 | 1.844 | 0.0100 | 9.06 | 10.0 | -9.4 | 30.0 |
| Hexachloroethane | Ave | 0.6881 | 0.6347 | 0.3000 | 9.22 | 10.0 | -7.8 | 30.0 |
| Nitrobenzene | Ave | 0.3728 | 0.3466 | 0.2000 | 9.30 | 10.0 | -7.0 | 30.0 |
| Isophorone | Ave | 0.6715 | 0.6521 | 0.4000 | 9.71 | 10.0 | -2.9 | 30.0 |
| 2-Nitrophenol | Ave | 0.1801 | 0.1714 | 0.1000 | 9.52 | 10.0 | -4.8 | 30.0 |
| 2,4-Dimethylphenol | Ave | 0.3720 | 0.3420 | 0.2000 | 9.19 | 10.0 | -8.1 | 30.0 |
| Benzoic acid | Ave | 0.1741 | 0.1903 | | 21.9 | 20.0 | 9.3 | 30.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4019 | 0.3824 | 0.3000 | 9.52 | 10.0 | -4.8 | 30.0 |
| 2,4-Dichlorophenol | Ave | 0.2928 | 0.2786 | 0.2000 | 9.51 | 10.0 | -4.9 | 30.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.3042 | 0.2929 | | 9.63 | 10.0 | -3.7 | 30.0 |
| Naphthalene | Ave | 1.048 | 0.995 | 0.7000 | 9.49 | 10.0 | -5.1 | 30.0 |
| 4-Chloroaniline | Ave | 0.4368 | 0.3976 | 0.0100 | 9.10 | 10.0 | -9.0 | 30.0 |
| 2,6-Dichlorophenol | Ave | 0.2922 | 0.2789 | | 9.55 | 10.0 | -4.5 | 30.0 |
| Hexachlorobutadiene | Ave | 0.1998 | 0.1851 | 0.0100 | 9.26 | 10.0 | -7.4 | 30.0 |
| Caprolactam | Ave | 0.1035 | 0.0982 | 0.0100 | 19.0 | 20.0 | -5.2 | 30.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3075 | 0.2976 | 0.2000 | 9.68 | 10.0 | -3.2 | 30.0 |
| 2-Methylnaphthalene | Ave | 0.7041 | 0.6346 | 0.4000 | 9.01 | 10.0 | -9.9 | 30.0 |
| 1-Methylnaphthalene | Ave | 0.6096 | 0.5922 | | 9.71 | 10.0 | -2.9 | 30.0 |
| Hexachlorocyclopentadiene | Ave | 0.3154 | 0.2915 | 0.0500 | 9.24 | 10.0 | -7.6 | 30.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.5061 | 0.4956 | 0.0100 | 9.79 | 10.0 | -2.1 | 30.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3644 | 0.3516 | 0.2000 | 9.65 | 10.0 | -3.5 | 30.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3740 | 0.3746 | 0.2000 | 10.0 | 10.0 | 0.2 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-206629/11 Calibration Date: 11/13/2015 13:12

Instrument ID: A4HP10 Calib Start Date: 11/13/2015 09:35

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/13/2015 13:36

Lab File ID: 51113011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.372 | 1.301 | 0.0100 | 9.48 | 10.0 | -5.2 | 30.0 |
| 2-Chloronaphthalene | Ave | 1.121 | 1.080 | 0.8000 | 9.64 | 10.0 | -3.6 | 30.0 |
| 2-Nitroaniline | Ave | 0.3783 | 0.3693 | 0.0100 | 9.76 | 10.0 | -2.4 | 30.0 |
| Dimethyl phthalate | Ave | 1.289 | 1.223 | 0.0100 | 9.49 | 10.0 | -5.1 | 30.0 |
| 1,3-Dinitrobenzene | Ave | 0.1899 | 0.1924 | | 10.1 | 10.0 | 1.3 | 30.0 |
| 2,6-Dinitrotoluene | Ave | 0.2830 | 0.2800 | | 9.89 | 10.0 | -1.1 | 30.0 |
| Acenaphthylene | Ave | 1.667 | 1.580 | 0.9000 | 9.48 | 10.0 | -5.2 | 30.0 |
| 3-Nitroaniline | Ave | 0.3032 | 0.2838 | 0.0100 | 9.36 | 10.0 | -6.4 | 30.0 |
| 2,4-Dinitrophenol | Qua | | 0.1079 | 0.0100 | 18.3 | 20.0 | -8.7 | 30.0 |
| Acenaphthene | Ave | 1.062 | 1.019 | 0.9000 | 9.59 | 10.0 | -4.1 | 30.0 |
| 4-Nitrophenol | Ave | 0.2092 | 0.1949 | | 18.6 | 20.0 | -6.8 | 30.0 |
| 2,4-Dinitrotoluene | Ave | 0.3688 | 0.3590 | 0.2000 | 9.73 | 10.0 | -2.7 | 30.0 |
| Dibenzofuran | Ave | 1.579 | 1.480 | 0.8000 | 9.38 | 10.0 | -6.2 | 30.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.3249 | 0.3152 | 0.0100 | 9.70 | 10.0 | -3.0 | 30.0 |
| Diethyl phthalate | Ave | 1.273 | 1.227 | 0.0100 | 9.64 | 10.0 | -3.6 | 30.0 |
| Hexadecane | Ave | 0.6559 | 0.6385 | | 9.73 | 10.0 | -2.7 | 30.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.5971 | 0.5547 | 0.4000 | 9.29 | 10.0 | -7.1 | 30.0 |
| 4-Nitroaniline | Ave | 0.3028 | 0.2847 | 0.0100 | 9.40 | 10.0 | -6.0 | 30.0 |
| Fluorene | Ave | 1.313 | 1.232 | 0.9000 | 9.38 | 10.0 | -6.2 | 30.0 |
| 4,6-Dinitro-2-methylphenol | Ave | 0.1107 | 0.1028 | 0.0100 | 18.6 | 20.0 | -7.1 | 30.0 |
| N-Nitrosodiphenylamine | Ave | 0.5996 | 0.5884 | 0.0100 | 19.6 | 20.0 | -1.9 | 30.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 0.8232 | 0.8297 | | 10.1 | 10.0 | 0.8 | 30.0 |
| Azobenzene | Ave | 0.8237 | 0.8298 | | 10.1 | 10.0 | 0.7 | 30.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.2290 | 0.2222 | 0.1000 | 9.70 | 10.0 | -3.0 | 30.0 |
| Hexachlorobenzene | Ave | 0.2551 | 0.2411 | 0.1000 | 9.45 | 10.0 | -5.5 | 30.0 |
| Atrazine | Ave | 0.2351 | 0.2134 | 0.0100 | 18.2 | 20.0 | -9.2 | 30.0 |
| Pentachlorophenol | Ave | 0.1678 | 0.1751 | 0.0500 | 20.9 | 20.0 | 4.3 | 30.0 |
| n-Octadecane | Ave | 1.038 | 0.9823 | | 9.46 | 10.0 | -5.4 | 30.0 |
| Phenanthrene | Ave | 1.147 | 1.055 | 0.7000 | 9.19 | 10.0 | -8.1 | 30.0 |
| Anthracene | Ave | 1.161 | 1.092 | 0.7000 | 9.41 | 10.0 | -5.9 | 30.0 |
| Carbazole | Ave | 1.066 | 0.9539 | 0.0100 | 8.95 | 10.0 | -10.5 | 30.0 |
| Di-n-butyl phthalate | Ave | 1.439 | 1.379 | 0.0100 | 9.58 | 10.0 | -4.2 | 30.0 |
| Fluoranthene | Ave | 1.253 | 1.187 | 0.6000 | 9.47 | 10.0 | -5.3 | 30.0 |
| Benzidine | Ave | 0.5133 | 0.4927 | | 19.2 | 20.0 | -4.0 | 30.0 |
| Pyrene | Ave | 1.123 | 1.039 | 0.6000 | 9.25 | 10.0 | -7.5 | 30.0 |
| Butyl benzyl phthalate | Ave | 0.5440 | 0.5398 | 0.0100 | 9.92 | 10.0 | -0.8 | 30.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3962 | 0.3861 | 0.0100 | 19.5 | 20.0 | -2.6 | 30.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.7789 | 0.7652 | 0.0100 | 9.82 | 10.0 | -1.8 | 30.0 |
| Benzo[a]anthracene | Ave | 1.081 | 1.022 | 0.8000 | 9.46 | 10.0 | -5.4 | 30.0 |
| Chrysene | Ave | 0.995 | 0.9698 | 0.7000 | 9.75 | 10.0 | -2.5 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-206629/11  Calibration Date: 11/13/2015 13:12

Instrument ID: A4HP10  Calib Start Date: 11/13/2015 09:35

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 11/13/2015 13:36

Lab File ID: 51113011.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Ave | 1.621 | 1.669 | 0.0100 | 10.3 | 10.0 | 3.0 | 30.0 |
| Benzo[b]fluoranthene | Ave | 1.177 | 1.172 | 0.7000 | 9.96 | 10.0 | -0.4 | 30.0 |
| Benzo[k]fluoranthene | Ave | 1.231 | 1.152 | 0.7000 | 9.36 | 10.0 | -6.4 | 30.0 |
| Benzo[a]pyrene | Ave | 1.130 | 1.092 | 0.7000 | 9.66 | 10.0 | -3.4 | 30.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.248 | 1.198 | 0.5000 | 9.60 | 10.0 | -4.0 | 30.0 |
| Dibenz(a,h)anthracene | Ave | 1.037 | 1.022 | 0.4000 | 9.85 | 10.0 | -1.5 | 30.0 |
| Benzo[g,h,i]perylene | Ave | 1.046 | 1.036 | 0.5000 | 9.90 | 10.0 | -1.0 | 30.0 |
| 2-Fluorophenol (Surr) | Ave | 1.361 | 1.380 | | 10.1 | 10.0 | 1.4 | 30.0 |
| Phenol-d5 (Surr) | Ave | 1.647 | 1.680 | | 10.2 | 10.0 | 2.0 | 30.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.4042 | 0.4213 | | 10.4 | 10.0 | 4.2 | 30.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.235 | 1.294 | | 10.5 | 10.0 | 4.8 | 30.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1856 | 0.1971 | | 10.6 | 10.0 | 6.2 | 30.0 |
| Terphenyl-d14 (Surr) | Ave | 0.8105 | 0.8171 | | 10.1 | 10.0 | 0.8 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207287/2 Calibration Date: 11/18/2015 08:26

Instrument ID: A4HP10 Calib Start Date: 11/13/2015 09:35

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/13/2015 13:36

Lab File ID: 51118002.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.5901 | 0.5726 | | 9.70 | 10.0 | -3.0 | 20.0 |
| N-Nitrosodimethylamine | Ave | 0.7688 | 0.7535 | | 9.80 | 10.0 | -2.0 | 20.0 |
| Pyridine | Ave | 1.381 | 1.381 | | 10.0 | 10.0 | -0.0 | 20.0 |
| Benzaldehyde | Ave | 1.142 | 1.103 | 0.0100 | 19.3 | 20.0 | -3.4 | 20.0 |
| Phenol | Ave | 1.754 | 1.726 | 0.8000 | 9.84 | 10.0 | -1.6 | 20.0 |
| Aniline | Ave | 2.152 | 2.181 | | 10.1 | 10.0 | 1.4 | 20.0 |
| Bis(2-chloroethyl)ether | Ave | 1.382 | 1.375 | 0.7000 | 9.95 | 10.0 | -0.5 | 20.0 |
| 2-Chlorophenol | Ave | 1.495 | 1.463 | 0.8000 | 9.79 | 10.0 | -2.1 | 20.0 |
| n-Decane | Ave | 2.087 | 1.998 | | 9.57 | 10.0 | -4.3 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.542 | 1.520 | | 9.86 | 10.0 | -1.4 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.578 | 1.567 | | 9.93 | 10.0 | -0.7 | 20.0 |
| Benzyl alcohol | Ave | 0.9548 | 0.9285 | | 9.72 | 10.0 | -2.8 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.511 | 1.482 | | 9.81 | 10.0 | -1.9 | 20.0 |
| 2-Methylphenol | Ave | 1.327 | 1.299 | 0.7000 | 9.78 | 10.0 | -2.2 | 20.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.522 | 2.325 | | 9.22 | 10.0 | -7.8 | 20.0 |
| Indene | Ave | 2.622 | 2.564 | | 19.6 | 20.0 | -2.2 | 20.0 |
| 3 & 4 Methylphenol | Ave | 1.390 | 1.376 | | 9.90 | 10.0 | -1.0 | 20.0 |
| N-Nitrosodi-n-propylamine | Ave | 0.9159 | 0.8578 | 0.5000 | 9.37 | 10.0 | -6.3 | 20.0 |
| Acetophenone | Ave | 2.035 | 1.962 | 0.0100 | 9.64 | 10.0 | -3.6 | 20.0 |
| Hexachloroethane | Ave | 0.6881 | 0.6762 | 0.3000 | 9.83 | 10.0 | -1.7 | 20.0 |
| Nitrobenzene | Ave | 0.3728 | 0.3700 | 0.2000 | 9.93 | 10.0 | -0.7 | 20.0 |
| Isophorone | Ave | 0.6715 | 0.6648 | 0.4000 | 9.90 | 10.0 | -1.0 | 20.0 |
| 2-Nitrophenol | Ave | 0.1801 | 0.1837 | 0.1000 | 10.2 | 10.0 | 2.0 | 20.0 |
| 2,4-Dimethylphenol | Ave | 0.3720 | 0.3674 | 0.2000 | 9.88 | 10.0 | -1.2 | 20.0 |
| Benzoic acid | Ave | 0.1741 | 0.1635 | | 18.8 | 20.0 | -6.1 | 20.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4019 | 0.3950 | 0.3000 | 9.83 | 10.0 | -1.7 | 20.0 |
| 2,4-Dichlorophenol | Ave | 0.2928 | 0.2990 | 0.2000 | 10.2 | 10.0 | 2.1 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.3042 | 0.3011 | | 9.90 | 10.0 | -1.0 | 20.0 |
| Naphthalene | Ave | 1.048 | 1.038 | 0.7000 | 9.90 | 10.0 | -1.0 | 20.0 |
| 4-Chloroaniline | Ave | 0.4368 | 0.4263 | 0.0100 | 9.76 | 10.0 | -2.4 | 20.0 |
| 2,6-Dichlorophenol | Ave | 0.2922 | 0.2906 | | 9.95 | 10.0 | -0.5 | 20.0 |
| Hexachlorobutadiene | Ave | 0.1998 | 0.1947 | 0.0100 | 9.74 | 10.0 | -2.6 | 20.0 |
| Caprolactam | Ave | 0.1035 | 0.1034 | 0.0100 | 20.0 | 20.0 | -0.1 | 20.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3075 | 0.3043 | 0.2000 | 9.90 | 10.0 | -1.0 | 20.0 |
| 2-Methylnaphthalene | Ave | 0.7041 | 0.6761 | 0.4000 | 9.60 | 10.0 | -4.0 | 20.0 |
| 1-Methylnaphthalene | Ave | 0.6096 | 0.5970 | | 9.79 | 10.0 | -2.1 | 20.0 |
| Hexachlorocyclopentadiene | Ave | 0.3154 | 0.3422 | 0.0500 | 10.9 | 10.0 | 8.5 | 20.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.5061 | 0.5005 | 0.0100 | 9.89 | 10.0 | -1.1 | 20.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3644 | 0.3488 | 0.2000 | 9.57 | 10.0 | -4.3 | 20.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3740 | 0.3662 | 0.2000 | 9.79 | 10.0 | -2.1 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207287/2 Calibration Date: 11/18/2015 08:26

Instrument ID: A4HP10 Calib Start Date: 11/13/2015 09:35

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/13/2015 13:36

Lab File ID: 51118002.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.372 | 1.312 | 0.0100 | 9.56 | 10.0 | -4.4 | 20.0 |
| 2-Chloronaphthalene | Ave | 1.121 | 1.065 | 0.8000 | 9.50 | 10.0 | -5.0 | 20.0 |
| 2-Nitroaniline | Ave | 0.3783 | 0.3729 | 0.0100 | 9.86 | 10.0 | -1.4 | 20.0 |
| Dimethyl phthalate | Ave | 1.289 | 1.250 | 0.0100 | 9.70 | 10.0 | -3.0 | 20.0 |
| 1,3-Dinitrobenzene | Ave | 0.1899 | 0.2002 | | 10.5 | 10.0 | 5.5 | 20.0 |
| 2,6-Dinitrotoluene | Ave | 0.2830 | 0.2745 | | 9.70 | 10.0 | -3.0 | 20.0 |
| Acenaphthylene | Ave | 1.667 | 1.659 | 0.9000 | 9.95 | 10.0 | -0.5 | 20.0 |
| 3-Nitroaniline | Ave | 0.3032 | 0.3043 | 0.0100 | 10.0 | 10.0 | 0.4 | 20.0 |
| 2,4-Dinitrophenol | Qua | | 0.1335 | 0.0100 | 21.7 | 20.0 | 8.4 | 20.0 |
| Acenaphthene | Ave | 1.062 | 1.023 | 0.9000 | 9.63 | 10.0 | -3.7 | 20.0 |
| 4-Nitrophenol | Ave | 0.2092 | 0.2027 | | 19.4 | 20.0 | -3.1 | 20.0 |
| 2,4-Dinitrotoluene | Ave | 0.3688 | 0.3671 | 0.2000 | 9.95 | 10.0 | -0.5 | 20.0 |
| Dibenzofuran | Ave | 1.579 | 1.528 | 0.8000 | 9.68 | 10.0 | -3.2 | 20.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.3249 | 0.3162 | 0.0100 | 9.73 | 10.0 | -2.7 | 20.0 |
| Diethyl phthalate | Ave | 1.273 | 1.213 | 0.0100 | 9.53 | 10.0 | -4.7 | 20.0 |
| Hexadecane | Ave | 0.6559 | 0.6269 | | 9.56 | 10.0 | -4.4 | 20.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.5971 | 0.5767 | 0.4000 | 9.66 | 10.0 | -3.4 | 20.0 |
| 4-Nitroaniline | Ave | 0.3028 | 0.3105 | 0.0100 | 10.3 | 10.0 | 2.5 | 20.0 |
| Fluorene | Ave | 1.313 | 1.259 | 0.9000 | 9.59 | 10.0 | -4.1 | 20.0 |
| 4,6-Dinitro-2-methylphenol | Ave | 0.1107 | 0.1240 | 0.0100 | 22.4 | 20.0 | 12.0 | 20.0 |
| N-Nitrosodiphenylamine | Ave | 0.5996 | 0.5882 | 0.0100 | 19.6 | 20.0 | -1.9 | 20.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 0.8232 | 0.8029 | | 9.75 | 10.0 | -2.5 | 20.0 |
| Azobenzene | Ave | 0.8237 | 0.8030 | | 9.75 | 10.0 | -2.5 | 20.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.2290 | 0.2289 | 0.1000 | 10.0 | 10.0 | -0.0 | 20.0 |
| Hexachlorobenzene | Ave | 0.2551 | 0.2411 | 0.1000 | 9.45 | 10.0 | -5.5 | 20.0 |
| Atrazine | Ave | 0.2351 | 0.2276 | 0.0100 | 19.4 | 20.0 | -3.2 | 20.0 |
| Pentachlorophenol | Ave | 0.1678 | 0.1647 | 0.0500 | 19.6 | 20.0 | -1.9 | 20.0 |
| n-Octadecane | Ave | 1.038 | 0.9844 | | 9.48 | 10.0 | -5.2 | 20.0 |
| Phenanthrene | Ave | 1.147 | 1.113 | 0.7000 | 9.70 | 10.0 | -3.0 | 20.0 |
| Anthracene | Ave | 1.161 | 1.140 | 0.7000 | 9.82 | 10.0 | -1.8 | 20.0 |
| Carbazole | Ave | 1.066 | 1.059 | 0.0100 | 9.93 | 10.0 | -0.7 | 20.0 |
| Di-n-butyl phthalate | Ave | 1.439 | 1.380 | 0.0100 | 9.59 | 10.0 | -4.1 | 20.0 |
| Fluoranthene | Ave | 1.253 | 1.232 | 0.6000 | 9.83 | 10.0 | -1.7 | 20.0 |
| Benzidine | Ave | 0.5133 | 0.5673 | | 22.1 | 20.0 | 10.5 | 20.0 |
| Pyrene | Ave | 1.123 | 1.091 | 0.6000 | 9.72 | 10.0 | -2.8 | 20.0 |
| Butyl benzyl phthalate | Ave | 0.5440 | 0.5266 | 0.0100 | 9.68 | 10.0 | -3.2 | 20.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3962 | 0.3966 | 0.0100 | 20.0 | 20.0 | 0.0 | 20.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.7789 | 0.7507 | 0.0100 | 9.64 | 10.0 | -3.6 | 20.0 |
| Benzo[a]anthracene | Ave | 1.081 | 1.045 | 0.8000 | 9.67 | 10.0 | -3.3 | 20.0 |
| Chrysene | Ave | 0.995 | 0.9561 | 0.7000 | 9.61 | 10.0 | -3.9 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207287/2 | Calibration Date: 11/18/2015 08:26

Instrument ID: A4HP10 | Calib Start Date: 11/13/2015 09:35

GC Column: DB-624 ID: 0.18(mm) | Calib End Date: 11/13/2015 13:36

Lab File ID: 51118002.D | Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Ave | 1.621 | 1.532 | 0.0100 | 9.45 | 10.0 | -5.5 | 20.0 |
| Benzo[b]fluoranthene | Ave | 1.177 | 1.184 | 0.7000 | 10.1 | 10.0 | 0.7 | 20.0 |
| Benzo[k]fluoranthene | Ave | 1.231 | 1.184 | 0.7000 | 9.62 | 10.0 | -3.8 | 20.0 |
| Benzo[a]pyrene | Ave | 1.130 | 1.126 | 0.7000 | 9.96 | 10.0 | -0.4 | 20.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.248 | 1.252 | 0.5000 | 10.0 | 10.0 | 0.3 | 20.0 |
| Dibenz(a,h)anthracene | Ave | 1.037 | 1.072 | 0.4000 | 10.3 | 10.0 | 3.4 | 20.0 |
| Benzo[g,h,i]perylene | Ave | 1.046 | 1.072 | 0.5000 | 10.2 | 10.0 | 2.5 | 20.0 |
| 2-Fluorophenol (Surr) | Ave | 1.361 | 1.369 | | 10.1 | 10.0 | 0.6 | 20.0 |
| Phenol-d5 (Surr) | Ave | 1.647 | 1.654 | | 10.0 | 10.0 | 0.4 | 20.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.4042 | 0.3889 | | 9.62 | 10.0 | -3.8 | 20.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.235 | 1.171 | | 9.49 | 10.0 | -5.1 | 20.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1856 | 0.1794 | | 9.67 | 10.0 | -3.3 | 20.0 |
| Terphenyl-d14 (Surr) | Ave | 0.8105 | 0.7725 | | 9.53 | 10.0 | -4.7 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207541/2 Calibration Date: 11/19/2015 09:50

Instrument ID: A4HP10 Calib Start Date: 11/13/2015 09:35

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/13/2015 13:36

Lab File ID: 51119002.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.5901 | 0.5882 | | 9.97 | 10.0 | -0.3 | 20.0 |
| N-Nitrosodimethylamine | Ave | 0.7688 | 0.7654 | | 9.96 | 10.0 | -0.4 | 20.0 |
| Pyridine | Ave | 1.381 | 1.421 | | 10.3 | 10.0 | 2.9 | 20.0 |
| Benzaldehyde | Ave | 1.142 | 1.113 | 0.0100 | 19.5 | 20.0 | -2.5 | 20.0 |
| Phenol | Ave | 1.754 | 1.764 | 0.8000 | 10.1 | 10.0 | 0.6 | 20.0 |
| Aniline | Ave | 2.152 | 2.156 | | 10.0 | 10.0 | 0.2 | 20.0 |
| Bis(2-chloroethyl)ether | Ave | 1.382 | 1.427 | 0.7000 | 10.3 | 10.0 | 3.3 | 20.0 |
| 2-Chlorophenol | Ave | 1.495 | 1.495 | 0.8000 | 10.0 | 10.0 | 0.0 | 20.0 |
| n-Decane | Ave | 2.087 | 2.036 | | 9.76 | 10.0 | -2.4 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.542 | 1.548 | | 10.0 | 10.0 | 0.4 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.578 | 1.568 | | 9.93 | 10.0 | -0.7 | 20.0 |
| Benzyl alcohol | Ave | 0.9548 | 0.9173 | | 9.61 | 10.0 | -3.9 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.511 | 1.505 | | 9.96 | 10.0 | -0.4 | 20.0 |
| 2-Methylphenol | Ave | 1.327 | 1.332 | 0.7000 | 10.0 | 10.0 | 0.3 | 20.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.522 | 2.340 | | 9.28 | 10.0 | -7.2 | 20.0 |
| Indene | Ave | 2.622 | 2.616 | | 20.0 | 20.0 | -0.2 | 20.0 |
| 3 & 4 Methylphenol | Ave | 1.390 | 1.403 | | 10.1 | 10.0 | 0.9 | 20.0 |
| N-Nitrosodi-n-propylamine | Ave | 0.9159 | 0.8920 | 0.5000 | 9.74 | 10.0 | -2.6 | 20.0 |
| Acetophenone | Ave | 2.035 | 2.022 | 0.0100 | 9.94 | 10.0 | -0.6 | 20.0 |
| Hexachloroethane | Ave | 0.6881 | 0.6614 | 0.3000 | 9.61 | 10.0 | -3.9 | 20.0 |
| Nitrobenzene | Ave | 0.3728 | 0.3624 | 0.2000 | 9.72 | 10.0 | -2.8 | 20.0 |
| Isophorone | Ave | 0.6715 | 0.6755 | 0.4000 | 10.1 | 10.0 | 0.6 | 20.0 |
| 2-Nitrophenol | Ave | 0.1801 | 0.1857 | 0.1000 | 10.3 | 10.0 | 3.1 | 20.0 |
| 2,4-Dimethylphenol | Ave | 0.3720 | 0.3589 | 0.2000 | 9.65 | 10.0 | -3.5 | 20.0 |
| Benzoic acid | Ave | 0.1741 | 0.1761 | | 20.2 | 20.0 | 1.2 | 20.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4019 | 0.4102 | 0.3000 | 10.2 | 10.0 | 2.1 | 20.0 |
| 2,4-Dichlorophenol | Ave | 0.2928 | 0.2957 | 0.2000 | 10.1 | 10.0 | 1.0 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.3042 | 0.3068 | | 10.1 | 10.0 | 0.8 | 20.0 |
| Naphthalene | Ave | 1.048 | 1.032 | 0.7000 | 9.84 | 10.0 | -1.6 | 20.0 |
| 4-Chloroaniline | Ave | 0.4368 | 0.4345 | 0.0100 | 9.95 | 10.0 | -0.5 | 20.0 |
| 2,6-Dichlorophenol | Ave | 0.2922 | 0.2926 | | 10.0 | 10.0 | 0.2 | 20.0 |
| Hexachlorobutadiene | Ave | 0.1998 | 0.1932 | 0.0100 | 9.67 | 10.0 | -3.3 | 20.0 |
| Caprolactam | Ave | 0.1035 | 0.1060 | 0.0100 | 20.5 | 20.0 | 2.3 | 20.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3075 | 0.3057 | 0.2000 | 9.94 | 10.0 | -0.6 | 20.0 |
| 2-Methylnaphthalene | Ave | 0.7041 | 0.6947 | 0.4000 | 9.87 | 10.0 | -1.3 | 20.0 |
| 1-Methylnaphthalene | Ave | 0.6096 | 0.6073 | | 9.96 | 10.0 | -0.4 | 20.0 |
| Hexachlorocyclopentadiene | Ave | 0.3154 | 0.3065 | 0.0500 | 9.72 | 10.0 | -2.8 | 20.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.5061 | 0.5125 | 0.0100 | 10.1 | 10.0 | 1.3 | 20.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3644 | 0.3655 | 0.2000 | 10.0 | 10.0 | 0.3 | 20.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3740 | 0.3818 | 0.2000 | 10.2 | 10.0 | 2.1 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207541/2 Calibration Date: 11/19/2015 09:50

Instrument ID: A4HP10 Calib Start Date: 11/13/2015 09:35

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/13/2015 13:36

Lab File ID: 51119002.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.372 | 1.340 | 0.0100 | 9.77 | 10.0 | -2.3 | 20.0 |
| 2-Chloronaphthalene | Ave | 1.121 | 1.106 | 0.8000 | 9.87 | 10.0 | -1.3 | 20.0 |
| 2-Nitroaniline | Ave | 0.3783 | 0.3763 | 0.0100 | 9.95 | 10.0 | -0.5 | 20.0 |
| Dimethyl phthalate | Ave | 1.289 | 1.295 | 0.0100 | 10.0 | 10.0 | 0.5 | 20.0 |
| 1,3-Dinitrobenzene | Ave | 0.1899 | 0.1964 | | 10.3 | 10.0 | 3.5 | 20.0 |
| 2,6-Dinitrotoluene | Ave | 0.2830 | 0.2869 | | 10.1 | 10.0 | 1.4 | 20.0 |
| Acenaphthylene | Ave | 1.667 | 1.687 | 0.9000 | 10.1 | 10.0 | 1.2 | 20.0 |
| 3-Nitroaniline | Ave | 0.3032 | 0.3022 | 0.0100 | 9.97 | 10.0 | -0.3 | 20.0 |
| 2,4-Dinitrophenol | Qua | | 0.1209 | 0.0100 | 20.0 | 20.0 | 0.1 | 20.0 |
| Acenaphthene | Ave | 1.062 | 1.052 | 0.9000 | 9.91 | 10.0 | -0.9 | 20.0 |
| 4-Nitrophenol | Ave | 0.2092 | 0.1975 | | 18.9 | 20.0 | -5.6 | 20.0 |
| 2,4-Dinitrotoluene | Ave | 0.3688 | 0.3859 | 0.2000 | 10.5 | 10.0 | 4.6 | 20.0 |
| Dibenzofuran | Ave | 1.579 | 1.580 | 0.8000 | 10.0 | 10.0 | 0.1 | 20.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.3249 | 0.3237 | 0.0100 | 9.96 | 10.0 | -0.4 | 20.0 |
| Diethyl phthalate | Ave | 1.273 | 1.286 | 0.0100 | 10.1 | 10.0 | 1.1 | 20.0 |
| Hexadecane | Ave | 0.6559 | 0.6230 | | 9.50 | 10.0 | -5.0 | 20.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.5971 | 0.6034 | 0.4000 | 10.1 | 10.0 | 1.1 | 20.0 |
| 4-Nitroaniline | Ave | 0.3028 | 0.3153 | 0.0100 | 10.4 | 10.0 | 4.1 | 20.0 |
| Fluorene | Ave | 1.313 | 1.316 | 0.9000 | 10.0 | 10.0 | 0.2 | 20.0 |
| 4,6-Dinitro-2-methylphenol | Ave | 0.1107 | 0.1106 | 0.0100 | 20.0 | 20.0 | -0.0 | 20.0 |
| N-Nitrosodiphenylamine | Ave | 0.5996 | 0.5952 | 0.0100 | 19.9 | 20.0 | -0.7 | 20.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 0.8232 | 0.8009 | | 9.72 | 10.0 | -2.7 | 20.0 |
| Azobenzene | Ave | 0.8237 | 0.8017 | | 9.73 | 10.0 | -2.7 | 20.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.2290 | 0.2321 | 0.1000 | 10.1 | 10.0 | 1.3 | 20.0 |
| Hexachlorobenzene | Ave | 0.2551 | 0.2502 | 0.1000 | 9.81 | 10.0 | -1.9 | 20.0 |
| Atrazine | Ave | 0.2351 | 0.2364 | 0.0100 | 20.1 | 20.0 | 0.5 | 20.0 |
| Pentachlorophenol | Ave | 0.1678 | 0.1624 | 0.0500 | 19.4 | 20.0 | -3.2 | 20.0 |
| n-Octadecane | Ave | 1.038 | 1.027 | | 9.90 | 10.0 | -1.0 | 20.0 |
| Phenanthrene | Ave | 1.147 | 1.115 | 0.7000 | 9.72 | 10.0 | -2.8 | 20.0 |
| Anthracene | Ave | 1.161 | 1.132 | 0.7000 | 9.76 | 10.0 | -2.4 | 20.0 |
| Carbazole | Ave | 1.066 | 1.069 | 0.0100 | 10.0 | 10.0 | 0.2 | 20.0 |
| Di-n-butyl phthalate | Ave | 1.439 | 1.416 | 0.0100 | 9.84 | 10.0 | -1.6 | 20.0 |
| Fluoranthene | Ave | 1.253 | 1.236 | 0.6000 | 9.87 | 10.0 | -1.3 | 20.0 |
| Benzidine | Ave | 0.5133 | 0.5352 | | 20.9 | 20.0 | 4.3 | 20.0 |
| Pyrene | Ave | 1.123 | 1.098 | 0.6000 | 9.78 | 10.0 | -2.2 | 20.0 |
| Butyl benzyl phthalate | Ave | 0.5440 | 0.5500 | 0.0100 | 10.1 | 10.0 | 1.1 | 20.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3962 | 0.3948 | 0.0100 | 19.9 | 20.0 | -0.4 | 20.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.7789 | 0.7805 | 0.0100 | 10.0 | 10.0 | 0.2 | 20.0 |
| Benzo[a]anthracene | Ave | 1.081 | 1.062 | 0.8000 | 9.83 | 10.0 | -1.7 | 20.0 |
| Chrysene | Ave | 0.995 | 0.9889 | 0.7000 | 9.94 | 10.0 | -0.6 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207541/2 Calibration Date: 11/19/2015 09:50

Instrument ID: A4HP10 Calib Start Date: 11/13/2015 09:35

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 11/13/2015 13:36

Lab File ID: 51119002.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Ave | 1.621 | 1.642 | 0.0100 | 10.1 | 10.0 | 1.3 | 20.0 |
| Benzo[b]fluoranthene | Ave | 1.177 | 1.193 | 0.7000 | 10.1 | 10.0 | 1.4 | 20.0 |
| Benzo[k]fluoranthene | Ave | 1.231 | 1.261 | 0.7000 | 10.2 | 10.0 | 2.5 | 20.0 |
| Benzo[a]pyrene | Ave | 1.130 | 1.165 | 0.7000 | 10.3 | 10.0 | 3.1 | 20.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.248 | 1.281 | 0.5000 | 10.3 | 10.0 | 2.6 | 20.0 |
| Dibenz(a,h)anthracene | Ave | 1.037 | 1.062 | 0.4000 | 10.2 | 10.0 | 2.4 | 20.0 |
| Benzo[g,h,i]perylene | Ave | 1.046 | 1.082 | 0.5000 | 10.3 | 10.0 | 3.4 | 20.0 |
| 2-Fluorophenol (Surr) | Ave | 1.361 | 1.362 | | 10.0 | 10.0 | 0.1 | 20.0 |
| Phenol-d5 (Surr) | Ave | 1.647 | 1.654 | | 10.0 | 10.0 | 0.4 | 20.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.4042 | 0.3851 | | 9.53 | 10.0 | -4.7 | 20.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.235 | 1.220 | | 9.88 | 10.0 | -1.2 | 20.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1856 | 0.1869 | | 10.1 | 10.0 | 0.7 | 20.0 |
| Terphenyl-d14 (Surr) | Ave | 0.8105 | 0.7851 | | 9.69 | 10.0 | -3.1 | 20.0 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206887/23-A

Matrix: Solid | Lab File ID: 51119007.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 11:50

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 3.5 | U | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 9.5 | U | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 25 | U | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 8.9 | U | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 20 | U | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 20 | U | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 21 | U | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 17 | U | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 21 | U | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 0.45 | U | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 8.2 | U | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 0.50 | U | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 11 | U | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 9.1 | U | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 8.3 | U | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 18 | U | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 16 | U | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 9.2 | U | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 13 | U | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 21 | U | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 17 | U | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 13 | U | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 26 | U | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 17 | U | 330 | 17 |
| 83-32-9 | Acenaphthene | 0.76 | U | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 0.35 | U | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 9.2 | U | 100 | 9.2 |
| 120-12-7 | Anthracene | 0.78 | U | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 9.1 | U | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 12 | U | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 0.63 | U | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 0.64 | U | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 0.59 | U | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.35 | U | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206887/23-A

Matrix: Solid | Lab File ID: 51119007.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 11:50

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.68 | U | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 22 | U | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.0 | U | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 34.7 | J | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 10 | U | 70 | 10 |
| 105-60-2 | Caprolactam | 37 | U | 330 | 37 |
| 86-74-8 | Carbazole | 27 | U | 50 | 27 |
| 218-01-9 | Chrysene | 1.1 | U | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.66 | U | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 0.66 | U | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 16 | U | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 17 | U | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 15 | U | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 7.9 | U | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 0.55 | U | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 0.53 | U | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 2.1 | U | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 5.6 | U | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 8.1 | U | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 9.0 | U | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.35 | U | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 13 | U | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 6.3 | U | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 21 | U | 50 | 21 |
| 91-20-3 | Naphthalene | 0.82 | U | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 2.2 | U | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 9.1 | U | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 0.73 | U | 6.7 | 0.73 |
| 108-95-2 | Phenol | 7.3 | U | 50 | 7.3 |
| 129-00-0 | Pyrene | 0.44 | U | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 20 | U | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206887/23-A

Matrix: Solid | Lab File ID: 51119007.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 11:50

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 77 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 72 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 67 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 65 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 74 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 43 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206908/21-A

Matrix: Solid | Lab File ID: 51118003.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30(g) | Date Analyzed: 11/18/2015 08:51

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207287 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 3.5 | U | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 9.5 | U | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 25 | U | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 8.9 | U | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 20 | U | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 20 | U | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 21 | U | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 17 | U | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 21 | U | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 0.45 | U | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 8.2 | U | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 0.50 | U | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 11 | U | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 9.1 | U | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 8.3 | U | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 18 | U | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 16 | U | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 9.2 | U | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 13 | U | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 21 | U | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 17 | U | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 13 | U | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 26 | U | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 17 | U | 330 | 17 |
| 83-32-9 | Acenaphthene | 0.76 | U | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 0.35 | U | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 9.2 | U | 100 | 9.2 |
| 120-12-7 | Anthracene | 0.78 | U | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 9.1 | U | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 12 | U | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 0.63 | U | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 0.64 | U | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 0.59 | U | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.35 | U | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206908/21-A

Matrix: Solid | Lab File ID: 51118003.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30(g) | Date Analyzed: 11/18/2015 08:51

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207287 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.68 | U | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 22 | U | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.0 | U | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 61.2 | J | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 10 | U | 70 | 10 |
| 105-60-2 | Caprolactam | 37 | U | 330 | 37 |
| 86-74-8 | Carbazole | 27 | U | 50 | 27 |
| 218-01-9 | Chrysene | 1.1 | U | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.66 | U | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 0.66 | U | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 16 | U | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 17 | U | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 15 | U | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 7.9 | U | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 0.55 | U | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 0.53 | U | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 2.1 | U | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 5.6 | U | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 8.1 | U | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 9.0 | U | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.35 | U | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 13 | U | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 6.3 | U | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 21 | U | 50 | 21 |
| 91-20-3 | Naphthalene | 0.82 | U | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 2.2 | U | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 9.1 | U | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 0.73 | U | 6.7 | 0.73 |
| 108-95-2 | Phenol | 7.3 | U | 50 | 7.3 |
| 129-00-0 | Pyrene | 0.44 | U | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 20 | U | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206908/21-A

Matrix: Solid | Lab File ID: 51118003.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30(g) | Date Analyzed: 11/18/2015 08:51

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207287 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 74 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 73 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 67 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 70 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 70 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 30 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206887/24-A

Matrix: Solid | Lab File ID: 51119008.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 12:13

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 486 | | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 490 | | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 484 | | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 468 | | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 499 | | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 428 | | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 583 | | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 537 | | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 539 | | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 491 | | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 509 | | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 503 | | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 500 | | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 507 | | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 509 | | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 770 | | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 450 | | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 727 | | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 539 | | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 518 | | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 358 | | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 513 | | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 534 | | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 929 | | 330 | 17 |
| 83-32-9 | Acenaphthene | 487 | | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 498 | | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 491 | | 100 | 9.2 |
| 120-12-7 | Anthracene | 510 | | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 1060 | | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 1010 | | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 517 | | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 509 | | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 509 | | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 518 | | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206887/24-A

Matrix: Solid | Lab File ID: 51119008.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 12:13

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 530 | | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 526 | | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 515 | | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 607 | | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 545 | | 70 | 10 |
| 105-60-2 | Caprolactam | 1100 | | 330 | 37 |
| 86-74-8 | Carbazole | 530 | | 50 | 27 |
| 218-01-9 | Chrysene | 538 | | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 504 | | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 499 | | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 544 | | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 523 | | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 549 | | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 515 | | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 528 | | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 491 | | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 518 | | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 490 | | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 320 | J | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 473 | | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 515 | | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 527 | | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 513 | | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 1030 | | 50 | 21 |
| 91-20-3 | Naphthalene | 505 | | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 508 | | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 665 | | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 516 | | 6.7 | 0.73 |
| 108-95-2 | Phenol | 509 | | 50 | 7.3 |
| 129-00-0 | Pyrene | 507 | | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 501 | | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206887/24-A

Matrix: Solid | Lab File ID: 51119008.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 08:48

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 12:13

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 78 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 80 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 73 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 80 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 77 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 69 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206908/22-A

Matrix: Solid | Lab File ID: 51118004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30(g) | Date Analyzed: 11/18/2015 09:15

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207287 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 442 | | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 446 | | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 387 | | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 286 | | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 461 | | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 467 | | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 611 | | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 488 | | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 480 | | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 463 | | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 445 | | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 458 | | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 453 | | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 485 | | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 461 | | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 732 | | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 432 | | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 779 | | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 488 | | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 474 | | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 349 | | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 467 | | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 503 | | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 881 | | 330 | 17 |
| 83-32-9 | Acenaphthene | 464 | | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 462 | | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 438 | | 100 | 9.2 |
| 120-12-7 | Anthracene | 471 | | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 971 | | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 909 | | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 466 | | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 466 | | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 452 | | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 476 | | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206908/22-A

Matrix: Solid | Lab File ID: 51118004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30(g) | Date Analyzed: 11/18/2015 09:15

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207287 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 469 | | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 486 | | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 453 | | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 704 | | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 489 | | 70 | 10 |
| 105-60-2 | Caprolactam | 1050 | | 330 | 37 |
| 86-74-8 | Carbazole | 488 | | 50 | 27 |
| 218-01-9 | Chrysene | 495 | | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 469 | | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 464 | | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 490 | | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 479 | | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 489 | | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 445 | | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 473 | | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 454 | | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 477 | | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 447 | | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 397 | | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 418 | | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 477 | | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 478 | | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 451 | | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 946 | | 50 | 21 |
| 91-20-3 | Naphthalene | 462 | | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 467 | | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 669 | | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 471 | | 6.7 | 0.73 |
| 108-95-2 | Phenol | 456 | | 50 | 7.3 |
| 129-00-0 | Pyrene | 475 | | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 448 | | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206908/22-A

Matrix: Solid | Lab File ID: 51118004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30(g) | Date Analyzed: 11/18/2015 09:15

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207287 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 71 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 71 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 70 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 70 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 70 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 42 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 MS | Lab Sample ID: 240-57930-5 MS

Matrix: Solid | Lab File ID: 51119015.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.06(g) | Date Analyzed: 11/19/2015 15:00

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 547 | | 59 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 561 | | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 554 | | 180 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 543 | | 180 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 567 | | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 554 | | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 752 | | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 582 | | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 567 | | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 547 | | 59 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 587 | | 59 | 9.7 |
| 91-57-6 | 2-Methylnaphthalene | 564 | | 7.9 | 0.59 |
| 95-48-7 | 2-Methylphenol | 587 | | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 558 | | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 553 | | 59 | 9.8 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 856 | | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 491 | | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 856 | | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 597 | | 59 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 598 | | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 401 | | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 567 | | 59 | 15 |
| 100-01-6 | 4-Nitroaniline | 510 | | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 1030 | | 390 | 20 |
| 83-32-9 | Acenaphthene | 557 | | 7.9 | 0.90 |
| 208-96-8 | Acenaphthylene | 562 | | 7.9 | 0.41 |
| 98-86-2 | Acetophenone | 559 | | 120 | 11 |
| 120-12-7 | Anthracene | 552 | | 7.9 | 0.92 |
| 1912-24-9 | Atrazine | 1150 | | 240 | 11 |
| 100-52-7 | Benzaldehyde | 1160 | | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 567 | | 7.9 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 564 | | 7.9 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 560 | | 7.9 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 561 | | 7.9 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 MS | Lab Sample ID: 240-57930-5 MS

Matrix: Solid | Lab File ID: 51119015.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.06(g) | Date Analyzed: 11/19/2015 15:00

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 576 | | 7.9 | 0.80 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 605 | | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 593 | | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 613 | | 83 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 601 | | 83 | 12 |
| 105-60-2 | Caprolactam | 1300 | | 390 | 44 |
| 86-74-8 | Carbazole | 537 | | 59 | 32 |
| 218-01-9 | Chrysene | 586 | | 7.9 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 564 | | 7.9 | 0.78 |
| 132-64-9 | Dibenzofuran | 562 | | 59 | 0.78 |
| 84-66-2 | Diethyl phthalate | 591 | | 83 | 19 |
| 131-11-3 | Dimethyl phthalate | 575 | | 83 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 585 | | 83 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 570 | | 83 | 9.3 |
| 206-44-0 | Fluoranthene | 574 | | 7.9 | 0.65 |
| 86-73-7 | Fluorene | 553 | | 7.9 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 551 | | 7.9 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 543 | | 59 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 315 | J | 390 | 9.5 |
| 67-72-1 | Hexachloroethane | 519 | | 59 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 558 | | 7.9 | 0.41 |
| 78-59-1 | Isophorone | 596 | | 59 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 570 | | 59 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 1140 | | 59 | 25 |
| 91-20-3 | Naphthalene | 563 | | 7.9 | 0.97 |
| 98-95-3 | Nitrobenzene | 571 | | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 863 | | 180 | 11 |
| 85-01-8 | Phenanthrene | 560 | | 7.9 | 0.86 |
| 108-95-2 | Phenol | 583 | | 59 | 8.6 |
| 129-00-0 | Pyrene | 579 | | 7.9 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 578 | | 470 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 MS | Lab Sample ID: 240-57930-5 MS

Matrix: Solid | Lab File ID: 51119015.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 30.06(g) | Date Analyzed: 11/19/2015 15:00

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 76 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 78 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 73 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 78 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 74 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 72 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 MSD | Lab Sample ID: 240-57930-5 MSD

Matrix: Solid | Lab File ID: 51119016.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 29.87(g) | Date Analyzed: 11/19/2015 15:24

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 534 | | 59 | 4.2 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 534 | | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 553 | | 180 | 30 |
| 88-06-2 | 2,4,6-Trichlorophenol | 506 | | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 568 | | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 535 | | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 741 | | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 593 | | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 567 | | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 545 | | 59 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 546 | | 59 | 9.7 |
| 91-57-6 | 2-Methylnaphthalene | 568 | | 7.9 | 0.59 |
| 95-48-7 | 2-Methylphenol | 554 | | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 568 | | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 555 | | 59 | 9.8 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 899 | | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 483 | | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 875 | | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 593 | | 59 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 579 | | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 419 | | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 571 | | 59 | 15 |
| 100-01-6 | 4-Nitroaniline | 524 | | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 1020 | | 390 | 20 |
| 83-32-9 | Acenaphthene | 542 | | 7.9 | 0.90 |
| 208-96-8 | Acenaphthylene | 552 | | 7.9 | 0.42 |
| 98-86-2 | Acetophenone | 538 | | 120 | 11 |
| 120-12-7 | Anthracene | 564 | | 7.9 | 0.93 |
| 1912-24-9 | Atrazine | 1190 | | 240 | 11 |
| 100-52-7 | Benzaldehyde | 1100 | | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 580 | | 7.9 | 0.75 |
| 50-32-8 | Benzo[a]pyrene | 588 | | 7.9 | 0.76 |
| 205-99-2 | Benzo[b]fluoranthene | 575 | | 7.9 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 571 | | 7.9 | 0.42 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 MSD | Lab Sample ID: 240-57930-5 MSD

Matrix: Solid | Lab File ID: 51119016.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 29.87(g) | Date Analyzed: 11/19/2015 15:24

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 598 | | 7.9 | 0.81 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 594 | | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 564 | | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 627 | | 83 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 617 | | 83 | 12 |
| 105-60-2 | Caprolactam | 1320 | | 390 | 44 |
| 86-74-8 | Carbazole | 545 | | 59 | 32 |
| 218-01-9 | Chrysene | 608 | | 7.9 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 575 | | 7.9 | 0.78 |
| 132-64-9 | Dibenzofuran | 552 | | 59 | 0.78 |
| 84-66-2 | Diethyl phthalate | 590 | | 83 | 19 |
| 131-11-3 | Dimethyl phthalate | 576 | | 83 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 593 | | 83 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 592 | | 83 | 9.4 |
| 206-44-0 | Fluoranthene | 578 | | 7.9 | 0.65 |
| 86-73-7 | Fluorene | 545 | | 7.9 | 0.63 |
| 118-74-1 | Hexachlorobenzene | 562 | | 7.9 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 548 | | 59 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 238 | J | 390 | 9.6 |
| 67-72-1 | Hexachloroethane | 485 | | 59 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 579 | | 7.9 | 0.42 |
| 78-59-1 | Isophorone | 586 | | 59 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 540 | | 59 | 7.5 |
| 86-30-6 | N-Nitrosodiphenylamine | 1130 | | 59 | 25 |
| 91-20-3 | Naphthalene | 559 | | 7.9 | 0.97 |
| 98-95-3 | Nitrobenzene | 562 | | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 793 | | 180 | 11 |
| 85-01-8 | Phenanthrene | 567 | | 7.9 | 0.87 |
| 108-95-2 | Phenol | 554 | | 59 | 8.7 |
| 129-00-0 | Pyrene | 586 | | 7.9 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 544 | | 470 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 MSD | Lab Sample ID: 240-57930-5 MSD

Matrix: Solid | Lab File ID: 51119016.D

Analysis Method: 8270D | Date Collected: 11/13/2015 10:00

Extract. Method: 3540C | Date Extracted: 11/16/2015 09:20

Sample wt/vol: 29.87(g) | Date Analyzed: 11/19/2015 15:24

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207541 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 75 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 74 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 71 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 73 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 72 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 64 | | 10-110 |

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A4HP10 Start Date: 11/13/2015 09:10

Analysis Batch Number: 206629 End Date: 11/13/2015 19:38

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| DFTPP 240-206629/1 | | 11/13/2015 09:10 | 1 | 51113001.D | DB-624 0.18(mm) |
| STD5 240-206629/2 IC | | 11/13/2015 09:35 | 1 | 51113002.D | DB-624 0.18(mm) |
| STD4 240-206629/3 IC | | 11/13/2015 09:59 | 1 | 51113003.D | DB-624 0.18(mm) |
| STD3 240-206629/4 IC | | 11/13/2015 10:23 | 1 | 51113004.D | DB-624 0.18(mm) |
| STD2 240-206629/5 IC | | 11/13/2015 10:48 | 1 | 51113005.D | DB-624 0.18(mm) |
| STD1 240-206629/6 IC | | 11/13/2015 11:12 | 1 | 51113006.D | DB-624 0.18(mm) |
| STD8 240-206629/8 IC | | 11/13/2015 12:00 | 1 | 51113008.D | DB-624 0.18(mm) |
| STD7 240-206629/9 IC | | 11/13/2015 12:24 | 1 | 51113009.D | DB-624 0.18(mm) |
| STD6 240-206629/10 ICIS | | 11/13/2015 12:48 | 1 | 51113010.D | DB-624 0.18(mm) |
| ICV 240-206629/11 | | 11/13/2015 13:12 | 1 | 51113011.D | DB-624 0.18(mm) |
| STD9 240-206629/12 IC | | 11/13/2015 13:36 | 1 | 51113012.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/13/2015 14:01 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/13/2015 14:25 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/13/2015 14:49 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/13/2015 15:13 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/13/2015 15:37 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/13/2015 16:01 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/13/2015 19:38 | 4 | | DB-624 0.18(mm) |

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A4HP10 Start Date: 11/18/2015 08:03

Analysis Batch Number: 207287 End Date: 11/18/2015 19:18

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| DFTPP 240-207287/1 | | 11/18/2015 08:03 | 1 | 51118001.D | DB-624 0.18(mm) |
| CCV 240-207287/2 CCVIS | | 11/18/2015 08:26 | 1 | 51118002.D | DB-624 0.18(mm) |
| MB 240-206908/21-A | | 11/18/2015 08:51 | 1 | 51118003.D | DB-624 0.18(mm) |
| LCS 240-206908/22-A | | 11/18/2015 09:15 | 1 | 51118004.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 10:03 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 10:27 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 10:51 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 11:15 | 6.667 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 11:40 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 12:04 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 12:28 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 12:52 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 18:30 | 20 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 18:54 | 20 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/18/2015 19:18 | 20 | | DB-624 0.18(mm) |

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Instrument ID: A4HP10  Start Date: 11/19/2015 09:29

Analysis Batch Number: 207541  End Date: 11/19/2015 20:57

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| DFTPP 240-207541/1 | | 11/19/2015 09:29 | 1 | 51119001.D | DB-624 0.18(mm) |
| CCV 240-207541/2 CCVIS | | 11/19/2015 09:50 | 1 | 51119002.D | DB-624 0.18(mm) |
| ZZZZZ | | 11/19/2015 11:02 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 11/19/2015 11:26 | 1 | | DB-624 0.18(mm) |
| MB 240-206887/23-A | | 11/19/2015 11:50 | 1 | 51119007.D | DB-624 0.18(mm) |
| LCS 240-206887/24-A | | 11/19/2015 12:13 | 1 | 51119008.D | DB-624 0.18(mm) |
| 240-57930-5 | F-VB-UST-P-1 | 11/19/2015 14:37 | 1 | 51119014.D | DB-624 0.18(mm) |
| 240-57930-5 MS | F-VB-UST-P-1 MS | 11/19/2015 15:00 | 1 | 51119015.D | DB-624 0.18(mm) |
| 240-57930-5 MSD | F-VB-UST-P-1 MSD | 11/19/2015 15:24 | 1 | 51119016.D | DB-624 0.18(mm) |
| 240-57930-6 | F-VB-UST-P-2 | 11/19/2015 15:48 | 1 | 51119017.D | DB-624 0.18(mm) |
| 240-57930-7 | F-VB-UST-P-3 | 11/19/2015 16:12 | 1 | 51119018.D | DB-624 0.18(mm) |
| 240-57930-8 | F-VB-UST-P-4 | 11/19/2015 16:36 | 1 | 51119019.D | DB-624 0.18(mm) |
| 240-57930-9 | F-VB-UST-P-5 | 11/19/2015 17:00 | 1 | 51119020.D | DB-624 0.18(mm) |
| 240-57930-10 | F-VB-UST-P-6 | 11/19/2015 17:23 | 1 | 51119021.D | DB-624 0.18(mm) |
| 240-57930-2 | F-VS-UST-P-2 | 11/19/2015 19:23 | 1 | 51119026.D | DB-624 0.18(mm) |
| 240-57930-1 | F-VS-UST-P-1 | 11/19/2015 19:46 | 1 | 51119027.D | DB-624 0.18(mm) |
| 240-57930-11 | F-VS-UST-P-DUP | 11/19/2015 20:10 | 1 | 51119028.D | DB-624 0.18(mm) |
| 240-57930-4 | F-VS-UST-P-4 | 11/19/2015 20:34 | 1 | 51119029.D | DB-624 0.18(mm) |
| 240-57930-3 | F-VS-UST-P-3 | 11/19/2015 20:57 | 2.5 | 51119030.D | DB-624 0.18(mm) |

# GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 206887  Batch Start Date: 11/16/15 11:50  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/17/15 03:50

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exBENZALDEHYD 00031 | exBNASPIKE 00053 | exBNASURR W 00028 | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57930-D-6 | F-VB-UST-P-2 | 3540C, 8270D | T | 30.05 g | 2 mL | | | 1 mL | |
| 240-57930-D-7 | F-VB-UST-P-3 | 3540C, 8270D | T | 30.21 g | 2 mL | | | 1 mL | |
| 240-57930-D-8 | F-VB-UST-P-4 | 3540C, 8270D | T | 29.80 g | 2 mL | | | 1 mL | |
| 240-57930-D-9 | F-VB-UST-P-5 | 3540C, 8270D | T | 29.71 g | 2 mL | | | 1 mL | |
| 240-57930-D-10 | F-VB-UST-P-6 | 3540C, 8270D | T | 30.14 g | 2 mL | | | 1 mL | |
| 240-57930-D-11 | F-VS-UST-P-DUP | 3540C, 8270D | T | 30.46 g | 2 mL | | | 1 mL | |
| MB 240-206887/23 | | 3540C, 8270D | | 30 g | 2 mL | | | 1 mL | |
| LCS 240-206887/24 | | 3540C, 8270D | | 30 g | 2 mL | 1 mL | 1 mL | 1 mL | |

| Batch Notes | |
|---|---|
| Balance ID | 1339912 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2287348 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 206908  Batch Start Date: 11/16/15 11:50  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/17/15 03:50

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exBENZALDEHYD 00031 | exBNASPIKE 00053 | exBNASURR W 00028 | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57930-D-1 | F-VS-UST-P-1 | 3540C, 8270D | T | 29.52 g | 2 mL | | | 1 mL | |
| 240-57930-D-2 | F-VS-UST-P-2 | 3540C, 8270D | T | 29.77 g | 2 mL | | | 1 mL | |
| 240-57930-D-3 | F-VS-UST-P-3 | 3540C, 8270D | T | 30.26 g | 2 mL | | | 1 mL | |
| 240-57930-D-4 | F-VS-UST-P-4 | 3540C, 8270D | T | 30.32 g | 2 mL | | | 1 mL | |
| 240-57930-D-5 | F-VB-UST-P-1 | 3540C, 8270D | T | 30.24 g | 2 mL | | | 1 mL | |
| 240-57930-D-5 MS | F-VB-UST-P-1 | 3540C, 8270D | T | 30.06 g | 2 mL | 1 mL | 1 mL | 1 mL | |
| 240-57930-D-5 MSD | F-VB-UST-P-1 | 3540C, 8270D | T | 29.87 g | 2 mL | 1 mL | 1 mL | 1 mL | |
| MB 240-206908/21 | | 3540C, 8270D | | 30 g | 2 mL | | | 1 mL | |
| LCS 240-206908/22 | | 3540C, 8270D | | 30 g | 2 mL | 1 mL | 1 mL | 1 mL | |

| Batch Notes | |
|---|---|
| Balance ID | 1339912 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2287348 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8015C - GRO

## Nonhalogenated Organics using GC/FID - Modified (Gasoline Range Organics) by Method 8015C

FORM II
GC VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): RTX-VRX ID: 0.45(mm)

| Client Sample ID | Lab Sample ID | TFT1 # |
|---|---|---|
| F-VS-UST-P-1 | 240-57930-1 | 69 |
| F-VS-UST-P-2 | 240-57930-2 | 78 |
| F-VS-UST-P-3 | 240-57930-3 | 75 |
| F-VS-UST-P-4 | 240-57930-4 | 67 |
| F-VB-UST-P-1 | 240-57930-5 | 72 |
| F-VB-UST-P-2 | 240-57930-6 | 64 |
| F-VB-UST-P-3 | 240-57930-7 | 72 |
| F-VB-UST-P-4 | 240-57930-8 | 53 |
| F-VB-UST-P-5 | 240-57930-9 | 51 |
| F-VB-UST-P-6 | 240-57930-10 | 72 |
| F-VS-UST-P-DUP | 240-57930-11 | 70 |
| | MB 240-206878/1-A | 71 |
| | MB 240-207078/1-A | 79 |
| | MB 240-207274/1-A | 78 |
| | LCS 240-206878/2-A | 72 |
| | LCS 240-207078/2-A | 101 |
| | LCS 240-207274/2-A | 101 |
| F-VS-UST-P-DUP MS | 240-57930-11 MS | 87 |
| | 240-57774-C-8-J MS | 100 |
| F-VS-UST-P-DUP MSD | 240-57930-11 MSD | 73 |
| | 240-57774-C-8-K MSD | 100 |

| | QC LIMITS |
|---|---|
| TFT = Trifluorotoluene (Surr) | 40-139 |

# Column to be used to flag recovery values

FORM II 8015C

FORM III
GC VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: Y0111610.D

Lab ID: LCS 240-206878/2-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 651 | 81 | 74-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111706.D

Lab ID: LCS 240-207078/2-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 889 | 111 | 74-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111818.D

Lab ID: LCS 240-207274/2-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 836 | 104 | 74-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: Y0111623.D

Lab ID: 240-57930-11 MS Client ID: F-VS-UST-P-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 10100 | 1800 | 9930 | 80 | 10-136 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111715.D

Lab ID: 240-57774-C-8-J MS Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 899 | 52 U | 906 | 101 | 10-136 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: Y0111624.D

Lab ID: 240-57930-11 MSD Client ID: F-VS-UST-P-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 9290 | 9080 | 78 | 9 | 40 | 10-136 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111716.D

Lab ID: 240-57774-C-8-K MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 896 | 902 | 101 | 0 | 40 | 10-136 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM IV
GC VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: MB 240-206878/1-A

Matrix: Solid  Date Extracted: 11/16/2015 08:12

Lab File ID:(1) Y0111609.D  Lab File ID:(2)

Date Analyzed:(1) 11/16/2015 10:10  Date Analyzed:(2)

Instrument ID:(1) YPID  Instrument ID:(2)

GC Column:(1) RTX-VRX  ID: 0.45(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-206878/2-A | 11/16/2015 10:45 | |
| F-VS-UST-P-1 | 240-57930-1 | 11/16/2015 11:18 | |
| F-VS-UST-P-4 | 240-57930-4 | 11/16/2015 13:01 | |
| F-VB-UST-P-1 | 240-57930-5 | 11/16/2015 13:36 | |
| F-VB-UST-P-2 | 240-57930-6 | 11/16/2015 14:12 | |
| F-VB-UST-P-3 | 240-57930-7 | 11/16/2015 15:19 | |
| F-VB-UST-P-4 | 240-57930-8 | 11/16/2015 15:56 | |
| F-VB-UST-P-5 | 240-57930-9 | 11/16/2015 16:30 | |
| F-VB-UST-P-6 | 240-57930-10 | 11/16/2015 17:05 | |
| F-VS-UST-P-DUP | 240-57930-11 | 11/16/2015 17:40 | |
| F-VS-UST-P-DUP MS | 240-57930-11 MS | 11/16/2015 18:14 | |
| F-VS-UST-P-DUP MSD | 240-57930-11 MSD | 11/16/2015 18:49 | |

FORM IV
GC VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: MB 240-207078/1-A

Matrix: Solid  Date Extracted: 11/17/2015 07:48

Lab File ID:(1) A0111705.D  Lab File ID:(2)

Date Analyzed:(1) 11/17/2015 09:24  Date Analyzed:(2)

Instrument ID:(1) AFID  Instrument ID:(2)

GC Column:(1) RTX-VRX  ID: 0.45(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-207078/2-A | 11/17/2015 10:01 | |
| F-VS-UST-P-2 | 240-57930-2 | 11/17/2015 10:39 | |
| | 240-57774-C-8-J MS | 11/17/2015 15:35 | |
| | 240-57774-C-8-K MSD | 11/17/2015 16:12 | |

FORM IV
GC VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: MB 240-207274/1-A

Matrix: Solid  Date Extracted: 11/18/2015 06:57

Lab File ID:(1) A0111817.D  Lab File ID:(2)

Date Analyzed:(1) 11/18/2015 15:48  Date Analyzed:(2)

Instrument ID:(1) AFID  Instrument ID:(2)

GC Column:(1) RTX-VRX  ID: 0.45(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-207274/2-A | 11/18/2015 16:26 | |
| F-VS-UST-P-3 | 240-57930-3 | 11/18/2015 17:04 | |

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: STD4 240-194508/10  Date Analyzed: 08/24/2015 15:53

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): A0082110.D  Heated Purge: (Y/N) N

Calibration ID: 30361

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| INITIAL CALIBRATION SURROGATE | | | | 10.34 | | |
| UPPER LIMIT | | | | 10.44 | | |
| LOWER LIMIT | | | | 10.24 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| STD4 240-194508/10 ICRT | | 08/24/2015 15:53 | A0082110.D | 10.34 | | |
| ICV 240-194508/13 | | 08/24/2015 17:45 | A0082113.D | 10.34 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCVRT 240-207037/4  Date Analyzed: 11/17/2015 07:24

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): A0111704.D  Heated Purge: (Y/N) N

Calibration ID: 30361

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 10.37 | | |
| UPPER LIMIT | | | | 10.47 | | |
| LOWER LIMIT | | | | 10.27 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCVRT 240-207037/4 | | 11/17/2015 07:24 | A0111704.D | 10.37 | | |
| MB 240-207078/1-A | | 11/17/2015 09:24 | A0111705.D | 10.36 | | |
| LCS 240-207078/2-A | | 11/17/2015 10:01 | A0111706.D | 10.37 | | |
| 240-57930-2 | F-VS-UST-P-2 | 11/17/2015 10:39 | A0111707.D | 10.37 | | |
| CCV 240-207037/11 | | 11/17/2015 13:07 | A0111711.D | 10.37 | | |
| 240-57774-C-8-J MS | | 11/17/2015 15:35 | A0111715.D | 10.37 | | |
| 240-57774-C-8-K MSD | | 11/17/2015 16:12 | A0111716.D | 10.37 | | |
| CCV 240-207037/17 | | 11/17/2015 16:49 | A0111717.D | 10.37 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCVRT 240-207246/4  Date Analyzed: 11/18/2015 07:31

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): A0111804.D  Heated Purge: (Y/N) N

Calibration ID: 30361

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 10.37 | | |
| UPPER LIMIT | | | | 10.47 | | |
| LOWER LIMIT | | | | 10.27 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCVRT 240-207246/4 | | 11/18/2015 07:31 | A0111804.D | 10.37 | | |
| CCV 240-207246/16 | | 11/18/2015 15:12 | A0111816.D | 10.37 | | |
| MB 240-207274/1-A | | 11/18/2015 15:48 | A0111817.D | 10.37 | | |
| LCS 240-207274/2-A | | 11/18/2015 16:26 | A0111818.D | 10.37 | | |
| 240-57930-3 | F-VS-UST-P-3 | 11/18/2015 17:04 | A0111819.D | 10.37 | | |
| CCV 240-207246/26 | | 11/18/2015 21:24 | A0111826.D | 10.37 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCVRT 240-206833/4  Date Analyzed: 11/16/2015 07:13

Instrument ID: YPID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): Y0111604.D  Heated Purge: (Y/N) N

Calibration ID: 31080

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 9.74 | | |
| UPPER LIMIT | | | | 9.84 | | |
| LOWER LIMIT | | | | 9.64 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCVRT 240-206833/4 | | 11/16/2015 07:13 | Y0111604.D | 9.74 | | |
| CCV 240-206833/8 | | 11/16/2015 09:37 | Y0111608.D | 9.75 | | |
| MB 240-206878/1-A | | 11/16/2015 10:10 | Y0111609.D | 9.73 | | |
| LCS 240-206878/2-A | | 11/16/2015 10:45 | Y0111610.D | 9.75 | | |
| 240-57930-1 | F-VS-UST-P-1 | 11/16/2015 11:18 | Y0111611.D | 9.73 | | |
| 240-57930-4 | F-VS-UST-P-4 | 11/16/2015 13:01 | Y0111614.D | 9.73 | | |
| 240-57930-5 | F-VB-UST-P-1 | 11/16/2015 13:36 | Y0111615.D | 9.72 | | |
| 240-57930-6 | F-VB-UST-P-2 | 11/16/2015 14:12 | Y0111616.D | 9.73 | | |
| CCV 240-206833/17 | | 11/16/2015 14:46 | Y0111617.D | 9.73 | | |
| 240-57930-7 | F-VB-UST-P-3 | 11/16/2015 15:19 | Y0111618.D | 9.73 | | |
| 240-57930-8 | F-VB-UST-P-4 | 11/16/2015 15:56 | Y0111619.D | 9.72 | | |
| 240-57930-9 | F-VB-UST-P-5 | 11/16/2015 16:30 | Y0111620.D | 9.72 | | |
| 240-57930-10 | F-VB-UST-P-6 | 11/16/2015 17:05 | Y0111621.D | 9.73 | | |
| 240-57930-11 | F-VS-UST-P-DUP | 11/16/2015 17:40 | Y0111622.D | 9.73 | | |
| 240-57930-11 MS | F-VS-UST-P-DUP MS | 11/16/2015 18:14 | Y0111623.D | 9.73 | | |
| 240-57930-11 MSD | F-VS-UST-P-DUP MSD | 11/16/2015 18:49 | Y0111624.D | 9.75 | | |
| CCV 240-206833/26 | | 11/16/2015 19:58 | Y0111626.D | 9.73 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-1 | Lab Sample ID: 240-57930-1

Matrix: Solid | Lab File ID: Y0111611.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:30

Sample wt/vol: 0.55(g) | Date Analyzed: 11/16/2015 11:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 3500 | | 1100 | 490 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 69 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: A0111707.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:35

Sample wt/vol: 0.48(g) | Date Analyzed: 11/17/2015 10:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207037 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 5500 | B | 1200 | 570 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-3 | Lab Sample ID: 240-57930-3

Matrix: Solid | Lab File ID: A0111819.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:40

Sample wt/vol: 5.02(g) | Date Analyzed: 11/18/2015 17:04

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 207246 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 110 | J | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 75 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-4 | Lab Sample ID: 240-57930-4

Matrix: Solid | Lab File ID: Y0111614.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:45

Sample wt/vol: 5.01(g) | Date Analyzed: 11/16/2015 13:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 16.9 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 67 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 | Lab Sample ID: 240-57930-5

Matrix: Solid | Lab File ID: Y0111615.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:00

Sample wt/vol: 5.05(g) | Date Analyzed: 11/16/2015 13:36

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.4 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 72 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-2 | Lab Sample ID: 240-57930-6

Matrix: Solid | Lab File ID: Y0111616.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:05

Sample wt/vol: 4.96(g) | Date Analyzed: 11/16/2015 14:12

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.5 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 64 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-3 | Lab Sample ID: 240-57930-7

Matrix: Solid | Lab File ID: Y0111618.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:10

Sample wt/vol: 5.08(g) | Date Analyzed: 11/16/2015 15:19

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 19.0 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 60 | J | 120 | 56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 72 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-4 | Lab Sample ID: 240-57930-8

Matrix: Solid | Lab File ID: Y0111619.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:15

Sample wt/vol: 4.95(g) | Date Analyzed: 11/16/2015 15:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 16.1 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 53 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-5 | Lab Sample ID: 240-57930-9

Matrix: Solid | Lab File ID: Y0111620.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:20

Sample wt/vol: 4.98(g) | Date Analyzed: 11/16/2015 16:30

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 16.7 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 51 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-6 | Lab Sample ID: 240-57930-10

Matrix: Solid | Lab File ID: Y0111621.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:25

Sample wt/vol: 4.96(g) | Date Analyzed: 11/16/2015 17:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 72 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP  Lab Sample ID: 240-57930-11

Matrix: Solid  Lab File ID: Y0111622.D

Analysis Method: 8015C  Date Collected: 11/13/2015 00:00

Sample wt/vol: 0.50(g)  Date Analyzed: 11/16/2015 17:40

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX  ID: 0.45(mm)

% Moisture: 15.6  Level: (low/med) Low

Analysis Batch No.: 206833  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 1800 | | 1200 | 540 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 70 | | 40-139 |

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 194508

SDG No.:

Instrument ID: AFID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 08/24/2015 13:59 Calibration End Date: 08/24/2015 17:07 Calibration ID: 30361

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-194508/7 | A0082107.D |
| Level 2 | STD2 240-194508/8 | A0082108.D |
| Level 3 | STD3 240-194508/9 | A0082109.D |
| Level 4 | STD4 240-194508/10 | A0082110.D |
| Level 5 | STD5 240-194508/11 | A0082111.D |
| Level 6 | STD6 240-194508/12 | A0082112.D |

| ANALYTE | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 6 | | | | | RT WINDOW | AVG RT |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 11.460 | 11.460 | 11.460 | 11.460 | 11.460 | 11.460 | | | | | 5.807 - 17.114 | 11.460 |
| C6-C12 | 12.725 | 12.725 | 12.725 | 12.725 | 12.725 | 12.725 | | | | | 5.807 - 19.644 | 12.725 |
| Trifluorotoluene (Surr) | 10.344 | 10.348 | 10.347 | 10.341 | 10.339 | 10.339 | | | | | 10.290 - 10.390 | 10.343 |

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 194508

SDG No.:

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 08/24/2015 13:59  Calibration End Date: 08/24/2015 17:07  Calibration ID: 30361

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-194508/7 | A0082107.D |
| Level 2 | STD2 240-194508/8 | A0082108.D |
| Level 3 | STD3 240-194508/9 | A0082109.D |
| Level 4 | STD4 240-194508/10 | A0082110.D |
| Level 5 | STD5 240-194508/11 | A0082111.D |
| Level 6 | STD6 240-194508/12 | A0082112.D |

| ANALYTE | CF | | | | CURVE TYPE | COEFFICIENT | | | # | MIN CF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 5 | LVL 2<br>LVL 6 | LVL 3 | LVL 4 | | B | M1 | M2 | | | | | | | | |
| Gasoline Range Organics [C6 - C10] | 281051<br>441376 | 404289<br>439409 | 445616 | 417800 | Ave | | 404923.648 | | | | 15.5 | | 20.0 | | | |
| C6-C12 | 344884<br>484706 | 442549<br>478304 | 494744 | 455788 | Ave | | 450162.318 | | | | 12.2 | | 20.0 | | | |
| Trifluorotoluene (Surr) | 452208<br>466474 | 509634<br>371094 | 436510 | 429984 | Ave | | 444317.252 | | | | 10.3 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 194508

SDG No.:

Instrument ID: AFID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 08/24/2015 13:59 Calibration End Date: 08/24/2015 17:07 Calibration ID: 30361

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-194508/7 | A0082107.D |
| Level 2 | STD2 240-194508/8 | A0082108.D |
| Level 3 | STD3 240-194508/9 | A0082109.D |
| Level 4 | STD4 240-194508/10 | A0082110.D |
| Level 5 | STD5 240-194508/11 | A0082111.D |
| Level 6 | STD6 240-194508/12 | A0082112.D |

| ANALYTE | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|
| | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Gasoline Range Organics [C6 - C10] | Ave | 28105105<br>703054501 | 80857808 | 178246588 | 334240056 | 529651436 | 100<br>1600 | 200 | 400 | 800 | 1200 |
| C6-C12 | Ave | 34488374<br>765287082 | 88509751 | 197897550 | 364630076 | 581646620 | 100<br>1600 | 200 | 400 | 800 | 1200 |
| Trifluorotoluene (Surr) | Ave | 452208<br>18554703 | 1019267 | 2182548 | 4299843 | 9329481 | 1.00<br>50.0 | 2.00 | 5.00 | 10.0 | 20.0 |

Curve Type Legend:
Ave = Average

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 201382

SDG No.:

Instrument ID: YPID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/12/2015 10:39 Calibration End Date: 10/12/2015 13:35 Calibration ID: 31080

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-201382/4 | Y0101204.D |
| Level 2 | STD2 240-201382/5 | Y0101205.D |
| Level 3 | STD3 240-201382/6 | Y0101206.D |
| Level 4 | STD4 240-201382/7 | Y0101207.D |
| Level 5 | STD5 240-201382/8 | Y0101208.D |
| Level 6 | STD6 240-201382/9 | Y0101209.D |

| ANALYTE | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 6 | | | | | RT WINDOW | AVG RT |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 12.032 | 12.032 | 12.032 | 12.032 | 12.032 | 12.032 | | | | | 4.928 - 19.136 | 12.032 |
| C6-C12 | 13.458 | 13.458 | 13.458 | 13.458 | 13.458 | 13.458 | | | | | 4.928 - 21.988 | 13.458 |
| Trifluorotoluene (Surr) | 9.720 | 9.722 | 9.720 | 9.726 | 9.719 | 9.722 | | | | | 9.476 - 9.976 | 9.722 |

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 201382

SDG No.:

Instrument ID: YPID  GC Column: RTX-VRX  ID: 0.45(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/12/2015 10:39  Calibration End Date: 10/12/2015 13:35  Calibration ID: 31080

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-201382/4 | Y0101204.D |
| Level 2 | STD2 240-201382/5 | Y0101205.D |
| Level 3 | STD3 240-201382/6 | Y0101206.D |
| Level 4 | STD4 240-201382/7 | Y0101207.D |
| Level 5 | STD5 240-201382/8 | Y0101208.D |
| Level 6 | STD6 240-201382/9 | Y0101209.D |

| ANALYTE | CF | | | | CURVE TYPE | COEFFICIENT | | | # | MIN CF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 5 | LVL 2<br>LVL 6 | LVL 3 | LVL 4 | | B | M1 | M2 | | | | | | | | |
| Gasoline Range Organics [C6 - C10] | 188416<br>173575 | 165916<br>171151 | 176757 | 171370 | Ave | | 174530.549 | | | | 4.4 | | 20.0 | | | |
| C6-C12 | 222173<br>187083 | 182916<br>184022 | 192254 | 185117 | Ave | | 192260.693 | | | | 7.8 | | 20.0 | | | |
| Trifluorotoluene (Surr) | 210718<br>163918 | 183500<br>159255 | 169226 | 148916 | Ave | | 172588.858 | | | | 12.7 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 201382

SDG No.:

Instrument ID: YPID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/12/2015 10:39 Calibration End Date: 10/12/2015 13:35 Calibration ID: 31080

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-201382/4 | Y0101204.D |
| Level 2 | STD2 240-201382/5 | Y0101205.D |
| Level 3 | STD3 240-201382/6 | Y0101206.D |
| Level 4 | STD4 240-201382/7 | Y0101207.D |
| Level 5 | STD5 240-201382/8 | Y0101208.D |
| Level 6 | STD6 240-201382/9 | Y0101209.D |

| ANALYTE | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|
| | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Gasoline Range Organics [C6 - C10] | Ave | 18841554<br>273840999 | 33183190 | 70702651 | 137095852 | 208289684 | 100<br>1600 | 200 | 400 | 800 | 1200 |
| C6-C12 | Ave | 22217288<br>294434655 | 36583184 | 76901726 | 148093223 | 224499426 | 100<br>1600 | 200 | 400 | 800 | 1200 |
| Trifluorotoluene (Surr) | Ave | 210718<br>7962770 | 367000 | 846131 | 1489159 | 3278353 | 1.00<br>50.0 | 2.00 | 5.00 | 10.0 | 20.0 |

Curve Type Legend:
Ave = Average

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-194508/13 Calibration Date: 08/24/2015 17:45

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0082113.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 425225 | | 840 | 800 | 5.0 | 20.0 |
| C6-C12 | Ave | 450162 | 466294 | | 829 | 800 | 3.6 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 487053 | | 11.0 | 10.0 | 9.6 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-194508/13 Calibration Date: 08/24/2015 17:45

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0082113.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.46 | 5.81 | 17.11 |
| C6-C12 | 12.73 | 5.81 | 19.64 |
| Trifluorotoluene (Surr) | 10.34 | 10.29 | 10.39 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVRT 240-207037/4 Calibration Date: 11/17/2015 07:24

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111704.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 446659 | | 882 | 800 | 10.3 | 20.0 |
| C6-C12 | Ave | 450162 | 487634 | | 867 | 800 | 8.3 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 421567 | | 9.49 | 10.0 | -5.1 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVRT 240-207037/4 Calibration Date: 11/17/2015 07:24

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111704.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.55 | 5.95 | 17.12 |
| C6-C12 | 12.79 | 5.95 | 19.67 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207037/11 Calibration Date: 11/17/2015 13:07

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111711.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 412790 | | 816 | 800 | 1.9 | 20.0 |
| C6-C12 | Ave | 450162 | 448627 | | 797 | 800 | -0.3 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 415089 | | 9.34 | 10.0 | -6.6 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207037/11 Calibration Date: 11/17/2015 13:07

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111711.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.55 | 5.95 | 17.12 |
| C6-C12 | 12.79 | 5.95 | 19.67 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207037/17 Calibration Date: 11/17/2015 16:49

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111717.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 402593 | | 795 | 800 | -0.6 | 20.0 |
| C6-C12 | Ave | 450162 | 437404 | | 777 | 800 | -2.8 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 399103 | | 8.98 | 10.0 | -10.2 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207037/17 Calibration Date: 11/17/2015 16:49

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111717.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.55 | 5.95 | 17.12 |
| C6-C12 | 12.79 | 5.95 | 19.67 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVRT 240-207246/4 Calibration Date: 11/18/2015 07:31

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111804.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 449468 | | 888 | 800 | 11.0 | 20.0 |
| C6-C12 | Ave | 450162 | 487761 | | 867 | 800 | 8.4 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 420667 | | 9.47 | 10.0 | -5.3 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVRT 240-207246/4 Calibration Date: 11/18/2015 07:31

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111804.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.55 | 5.94 | 17.18 |
| C6-C12 | 12.79 | 5.94 | 19.70 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207246/16 Calibration Date: 11/18/2015 15:12

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111816.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 411988 | | 814 | 800 | 1.7 | 20.0 |
| C6-C12 | Ave | 450162 | 448100 | | 796 | 800 | -0.5 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 407157 | | 9.16 | 10.0 | -8.4 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207246/16 Calibration Date: 11/18/2015 15:12

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111816.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.55 | 5.94 | 17.18 |
| C6-C12 | 12.79 | 5.94 | 19.70 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207246/26 Calibration Date: 11/18/2015 21:24

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111826.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 407247 | | 805 | 800 | 0.6 | 20.0 |
| C6-C12 | Ave | 450162 | 442060 | | 786 | 800 | -1.8 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 393963 | | 8.87 | 10.0 | -11.3 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207246/26 Calibration Date: 11/18/2015 21:24

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111826.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.56 | 5.94 | 17.18 |
| C6-C12 | 12.82 | 5.94 | 19.70 |
| Trifluorotoluene (Surr) | 10.37 | 10.32 | 10.42 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-201382/10 Calibration Date: 10/12/2015 14:09

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0101210.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 174531 | 167422 | | 767 | 800 | -4.1 | 20.0 |
| C6-C12 | Ave | 192261 | 181216 | | 754 | 800 | -5.7 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 172589 | 166182 | | 9.63 | 10.0 | -3.7 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-201382/10 Calibration Date: 10/12/2015 14:09

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0101210.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 12.03 | 4.93 | 19.14 |
| C6-C12 | 13.46 | 4.93 | 21.99 |
| Trifluorotoluene (Surr) | 9.72 | 9.48 | 9.98 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVRT 240-206833/4 Calibration Date: 11/16/2015 07:13

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0111604.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 174531 | 175632 | | 805 | 800 | 0.6 | 20.0 |
| C6-C12 | Ave | 192261 | 190523 | | 793 | 800 | -0.9 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 172589 | 144962 | | 8.40 | 10.0 | -16.0 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVRT 240-206833/4 Calibration Date: 11/16/2015 07:13

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0111604.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 12.05 | 4.97 | 18.88 |
| C6-C12 | 13.49 | 4.97 | 21.97 |
| Trifluorotoluene (Surr) | 9.74 | 9.49 | 9.99 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-206833/8 Calibration Date: 11/16/2015 09:37

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0111608.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 174531 | 145133 | | 665 | 800 | -16.8 | 20.0 |
| C6-C12 | Ave | 192261 | 154736 | | 644 | 800 | -19.5 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 172589 | 159579 | | 9.25 | 10.0 | -7.5 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-206833/8 Calibration Date: 11/16/2015 09:37

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0111608.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 12.05 | 4.97 | 18.88 |
| C6-C12 | 13.49 | 4.97 | 21.97 |
| Trifluorotoluene (Surr) | 9.75 | 9.47 | 9.97 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-206833/17 Calibration Date: 11/16/2015 14:46

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0111617.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 174531 | 161427 | | 740 | 800 | -7.5 | 20.0 |
| C6-C12 | Ave | 192261 | 173583 | | 722 | 800 | -9.7 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 172589 | 140302 | | 8.13 | 10.0 | -18.7 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-206833/17 Calibration Date: 11/16/2015 14:46

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0111617.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 12.05 | 4.97 | 18.88 |
| C6-C12 | 13.49 | 4.97 | 21.97 |
| Trifluorotoluene (Surr) | 9.73 | 9.47 | 9.97 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-206833/26 Calibration Date: 11/16/2015 19:58

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0111626.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 174531 | 152687 | | 700 | 800 | -12.5 | 20.0 |
| C6-C12 | Ave | 192261 | 165483 | | 689 | 800 | -13.9 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 172589 | 138180 | | 8.01 | 10.0 | -19.9 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-206833/26 Calibration Date: 11/16/2015 19:58

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0111626.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 11.92 | 4.97 | 18.88 |
| C6-C12 | 13.47 | 4.97 | 21.97 |
| Trifluorotoluene (Surr) | 9.73 | 9.47 | 9.97 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-206878/1-A

Matrix: Solid  Lab File ID: Y0111609.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/16/2015 10:10

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206833  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 46 | U | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 71 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-207078/1-A

Matrix: Solid  Lab File ID: A0111705.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/17/2015 09:24

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 207037  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 47.9 | J | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 79 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-207274/1-A

Matrix: Solid  Lab File ID: A0111817.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/18/2015 15:48

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 207246  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 46 | U | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-206878/2-A

Matrix: Solid  Lab File ID: Y0111610.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/16/2015 10:45

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206833  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 651 | | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 72 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-207078/2-A

Matrix: Solid  Lab File ID: A0111706.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/17/2015 10:01

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 207037  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 889 | | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 101 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-207274/2-A

Matrix: Solid  Lab File ID: A0111818.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/18/2015 16:26

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 207246  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 836 | | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 101 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP MS | Lab Sample ID: 240-57930-11 MS

Matrix: Solid | Lab File ID: Y0111623.D

Analysis Method: 8015C | Date Collected: 11/13/2015 00:00

Sample wt/vol: 0.47(g) | Date Analyzed: 11/16/2015 18:14

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 9930 | | 1300 | 580 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 87 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57774-C-8-J MS

Matrix: Solid | Lab File ID: A0111715.D

Analysis Method: 8015C | Date Collected: 11/10/2015 13:30

Sample wt/vol: 4.92(g) | Date Analyzed: 11/17/2015 15:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 9.6 | Level: (low/med) Low

Analysis Batch No.: 207037 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 906 | | 110 | 52 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 100 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP MSD | Lab Sample ID: 240-57930-11 MSD

Matrix: Solid | Lab File ID: Y0111624.D

Analysis Method: 8015C | Date Collected: 11/13/2015 00:00

Sample wt/vol: 0.51(g) | Date Analyzed: 11/16/2015 18:49

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206833 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 9080 | | 1200 | 530 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 73 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57774-C-8-K MSD

Matrix: Solid | Lab File ID: A0111716.D

Analysis Method: 8015C | Date Collected: 11/10/2015 13:30

Sample wt/vol: 4.94(g) | Date Analyzed: 11/17/2015 16:12

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 9.6 | Level: (low/med) Low

Analysis Batch No.: 207037 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 902 | | 110 | 52 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 100 | | 40-139 |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Instrument ID: AFID  Start Date: 08/24/2015 13:59

Analysis Batch Number: 194508  End Date: 08/24/2015 17:45

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-194508/7 IC | | 08/24/2015 13:59 | 1 | A0082107.D | RTX-VRX 0.45(mm) |
| STD2 240-194508/8 IC | | 08/24/2015 14:37 | 1 | A0082108.D | RTX-VRX 0.45(mm) |
| STD3 240-194508/9 IC | | 08/24/2015 15:14 | 1 | A0082109.D | RTX-VRX 0.45(mm) |
| STD4 240-194508/10 ICRT | | 08/24/2015 15:53 | 1 | A0082110.D | RTX-VRX 0.45(mm) |
| STD5 240-194508/11 IC | | 08/24/2015 16:31 | 1 | A0082111.D | RTX-VRX 0.45(mm) |
| STD6 240-194508/12 IC | | 08/24/2015 17:07 | 1 | A0082112.D | RTX-VRX 0.45(mm) |
| ICV 240-194508/13 | | 08/24/2015 17:45 | 1 | A0082113.D | RTX-VRX 0.45(mm) |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: AFID Start Date: 11/17/2015 07:24

Analysis Batch Number: 207037 End Date: 11/17/2015 16:49

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCVRT 240-207037/4 | | 11/17/2015 07:24 | 1 | A0111704.D | RTX-VRX 0.45(mm) |
| MB 240-207078/1-A | | 11/17/2015 09:24 | 1 | A0111705.D | RTX-VRX 0.45(mm) |
| LCS 240-207078/2-A | | 11/17/2015 10:01 | 1 | A0111706.D | RTX-VRX 0.45(mm) |
| 240-57930-2 | F-VS-UST-P-2 | 11/17/2015 10:39 | 1 | A0111707.D | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/17/2015 11:16 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/17/2015 11:53 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/17/2015 12:30 | 1 | | RTX-VRX 0.45(mm) |
| CCV 240-207037/11 | | 11/17/2015 13:07 | 1 | A0111711.D | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/17/2015 13:44 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/17/2015 14:21 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/17/2015 14:58 | 1 | | RTX-VRX 0.45(mm) |
| 240-57774-C-8-J MS | | 11/17/2015 15:35 | 1 | A0111715.D | RTX-VRX 0.45(mm) |
| 240-57774-C-8-K MSD | | 11/17/2015 16:12 | 1 | A0111716.D | RTX-VRX 0.45(mm) |
| CCV 240-207037/17 | | 11/17/2015 16:49 | 1 | A0111717.D | RTX-VRX 0.45(mm) |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: AFID Start Date: 11/18/2015 07:31

Analysis Batch Number: 207246 End Date: 11/18/2015 21:24

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCVRT 240-207246/4 | | 11/18/2015 07:31 | 1 | A0111804.D | RTX-VRX 0.45(mm) |
| CCV 240-207246/16 | | 11/18/2015 15:12 | 1 | A0111816.D | RTX-VRX 0.45(mm) |
| MB 240-207274/1-A | | 11/18/2015 15:48 | 1 | A0111817.D | RTX-VRX 0.45(mm) |
| LCS 240-207274/2-A | | 11/18/2015 16:26 | 1 | A0111818.D | RTX-VRX 0.45(mm) |
| 240-57930-3 | F-VS-UST-P-3 | 11/18/2015 17:04 | 1 | A0111819.D | RTX-VRX 0.45(mm) |
| CCV 240-207246/26 | | 11/18/2015 21:24 | 1 | A0111826.D | RTX-VRX 0.45(mm) |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: YPID Start Date: 10/12/2015 10:39

Analysis Batch Number: 201382 End Date: 10/12/2015 15:19

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-201382/4 IC | | 10/12/2015 10:39 | 1 | Y0101204.D | RTX-VRX 0.45(mm) |
| STD2 240-201382/5 IC | | 10/12/2015 11:13 | 1 | Y0101205.D | RTX-VRX 0.45(mm) |
| STD3 240-201382/6 IC | | 10/12/2015 11:49 | 1 | Y0101206.D | RTX-VRX 0.45(mm) |
| STD4 240-201382/7 IC | | 10/12/2015 12:23 | 1 | Y0101207.D | RTX-VRX 0.45(mm) |
| STD5 240-201382/8 IC | | 10/12/2015 12:59 | 1 | Y0101208.D | RTX-VRX 0.45(mm) |
| STD6 240-201382/9 IC | | 10/12/2015 13:35 | 1 | Y0101209.D | RTX-VRX 0.45(mm) |
| ICV 240-201382/10 | | 10/12/2015 14:09 | 1 | Y0101210.D | RTX-VRX 0.45(mm) |
| ZZZZZ | | 10/12/2015 14:45 | 2 | | RTX-VRX 0.45(mm) |
| CCV 240-201382/12 | | 10/12/2015 15:19 | 1 | | RTX-VRX 0.45(mm) |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: YPID Start Date: 11/16/2015 07:13

Analysis Batch Number: 206833 End Date: 11/16/2015 19:58

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCVRT 240-206833/4 | | 11/16/2015 07:13 | 1 | Y0111604.D | RTX-VRX 0.45(mm) |
| CCV 240-206833/8 | | 11/16/2015 09:37 | 1 | Y0111608.D | RTX-VRX 0.45(mm) |
| MB 240-206878/1-A | | 11/16/2015 10:10 | 1 | Y0111609.D | RTX-VRX 0.45(mm) |
| LCS 240-206878/2-A | | 11/16/2015 10:45 | 1 | Y0111610.D | RTX-VRX 0.45(mm) |
| 240-57930-1 | F-VS-UST-P-1 | 11/16/2015 11:18 | 1 | Y0111611.D | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/16/2015 11:53 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/16/2015 12:28 | 1 | | RTX-VRX 0.45(mm) |
| 240-57930-4 | F-VS-UST-P-4 | 11/16/2015 13:01 | 1 | Y0111614.D | RTX-VRX 0.45(mm) |
| 240-57930-5 | F-VB-UST-P-1 | 11/16/2015 13:36 | 1 | Y0111615.D | RTX-VRX 0.45(mm) |
| 240-57930-6 | F-VB-UST-P-2 | 11/16/2015 14:12 | 1 | Y0111616.D | RTX-VRX 0.45(mm) |
| CCV 240-206833/17 | | 11/16/2015 14:46 | 1 | Y0111617.D | RTX-VRX 0.45(mm) |
| 240-57930-7 | F-VB-UST-P-3 | 11/16/2015 15:19 | 1 | Y0111618.D | RTX-VRX 0.45(mm) |
| 240-57930-8 | F-VB-UST-P-4 | 11/16/2015 15:56 | 1 | Y0111619.D | RTX-VRX 0.45(mm) |
| 240-57930-9 | F-VB-UST-P-5 | 11/16/2015 16:30 | 1 | Y0111620.D | RTX-VRX 0.45(mm) |
| 240-57930-10 | F-VB-UST-P-6 | 11/16/2015 17:05 | 1 | Y0111621.D | RTX-VRX 0.45(mm) |
| 240-57930-11 | F-VS-UST-P-DUP | 11/16/2015 17:40 | 1 | Y0111622.D | RTX-VRX 0.45(mm) |
| 240-57930-11 MS | F-VS-UST-P-DUP MS | 11/16/2015 18:14 | 1 | Y0111623.D | RTX-VRX 0.45(mm) |
| 240-57930-11 MSD | F-VS-UST-P-DUP MSD | 11/16/2015 18:49 | 1 | Y0111624.D | RTX-VRX 0.45(mm) |
| CCV 240-206833/26 | | 11/16/2015 19:58 | 1 | Y0111626.D | RTX-VRX 0.45(mm) |

# GC VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 206878  Batch Start Date: 11/16/15 08:12  Batch Analyst: Reed, Rachel T

Batch Method: 5030A  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| MB 240-206878/1 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| LCS 240-206878/2 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| 240-57930-C-1 | F-VS-UST-P-1 | 5030A, 8015C | T | 0.55 g | 5 mL | | | | |
| 240-57930-C-4 | F-VS-UST-P-4 | 5030A, 8015C | T | 5.01 g | 5 mL | | | | |
| 240-57930-C-5 | F-VB-UST-P-1 | 5030A, 8015C | T | 5.05 g | 5 mL | | | | |
| 240-57930-C-6 | F-VB-UST-P-2 | 5030A, 8015C | T | 4.96 g | 5 mL | | | | |
| 240-57930-C-7 | F-VB-UST-P-3 | 5030A, 8015C | T | 5.08 g | 5 mL | | | | |
| 240-57930-C-8 | F-VB-UST-P-4 | 5030A, 8015C | T | 4.95 g | 5 mL | | | | |
| 240-57930-C-9 | F-VB-UST-P-5 | 5030A, 8015C | T | 4.98 g | 5 mL | | | | |
| 240-57930-C-10 | F-VB-UST-P-6 | 5030A, 8015C | T | 4.96 g | 5 mL | | | | |
| 240-57930-C-11 | F-VS-UST-P-DUP | 5030A, 8015C | T | 0.50 g | 5 mL | | | | |
| 240-57930-C-11 MS | F-VS-UST-P-DUP | 5030A, 8015C | T | 0.47 g | 5 mL | | | | |
| 240-57930-C-11 MSD | F-VS-UST-P-DUP | 5030A, 8015C | T | 0.51 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| Balance ID | B045 |
| Blank Soil Lot Number | 1757979 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 207078  Batch Start Date: 11/17/15 07:48  Batch Analyst: Reed, Rachel T

Batch Method: 5030A  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| MB 240-207078/1 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| LCS 240-207078/2 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| 240-57930-C-2 | F-VS-UST-P-2 | 5030A, 8015C | T | 0.48 g | 5 mL | | | | |
| 240-57774-C-8 MS | | 5030A, 8015C | T | 4.92 g | 5 mL | | | | |
| 240-57774-C-8 MSD | | 5030A, 8015C | T | 4.94 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| Balance ID | B045 |
| Blank Soil Lot Number | 1757979 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 207274  Batch Start Date: 11/18/15 06:57  Batch Analyst: Reed, Rachel T

Batch Method: 5030A  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| MB 240-207274/1 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| LCS 240-207274/2 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| 240-57930-C-3 | F-VS-UST-P-3 | 5030A, 8015C | T | 5.02 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| Balance ID | B045 |
| Blank Soil Lot Number | 1757979 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8015C - DRO

## Nonhalogenated Organics using GC/FID - Modified (Diesel Range Organics) by Method 8015C

FORM II
DIESEL RANGE ORGANICS SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): ID:

| Client Sample ID | Lab Sample ID | OTPH1 # |
|---|---|---|
| F-VS-UST-P-1 | 240-57930-1 | 82 |
| F-VS-UST-P-2 | 240-57930-2 | 95 |
| F-VS-UST-P-3 | 240-57930-3 | 75 |
| F-VS-UST-P-4 | 240-57930-4 | 82 |
| F-VB-UST-P-1 | 240-57930-5 | 76 |
| F-VB-UST-P-2 | 240-57930-6 | 79 |
| F-VB-UST-P-3 | 240-57930-7 | 69 |
| F-VB-UST-P-4 | 240-57930-8 | 90 |
| F-VB-UST-P-5 | 240-57930-9 | 81 |
| F-VB-UST-P-6 | 240-57930-10 | 83 |
| F-VS-UST-P-DUP | 240-57930-11 | 79 |
| | MB 240-206884/23-A | 77 |
| | MB 240-207121/7-A | 75 |
| | LCS 240-206884/24-A | 82 |
| | LCS 240-207121/8-A | 103 |
| F-VS-UST-P-DUP MS | 240-57930-11 MS | 93 |
| F-VS-UST-P-DUP MSD | 240-57930-11 MSD | 93 |

| | QC LIMITS |
|---|---|
| OTPH = o-Terphenyl (Surr) | 40-160 |

# Column to be used to flag recovery values

FORM II 8015C

FORM III
DIESEL RANGE ORGANICS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11709.D

Lab ID: LCS 240-206884/24-A Client ID:

| COMPOUND | SPIKE ADDED (mg/Kg) | LCS CONCENTRATION (mg/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Diesel | 84.0 | 66.9 | 80 | 52-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
DIESEL RANGE ORGANICS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11816.D

Lab ID: LCS 240-207121/8-A Client ID:

| COMPOUND | SPIKE ADDED (mg/Kg) | LCS CONCENTRATION (mg/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Diesel | 84.0 | 83.2 | 99 | 52-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
DIESEL RANGE ORGANICS MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11821.D

Lab ID: 240-57930-11 MS Client ID: F-VS-UST-P-DUP MS

| COMPOUND | SPIKE ADDED (mg/Kg) | SAMPLE CONCENTRATION (mg/Kg) | MS CONCENTRATION (mg/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Diesel | 99.9 | 34 | 108 | 74 | 10-124 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
DIESEL RANGE ORGANICS MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11822.D

Lab ID: 240-57930-11 MSD Client ID: F-VS-UST-P-DUP MSD

| COMPOUND | SPIKE ADDED (mg/Kg) | MSD CONCENTRATION (mg/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Diesel | 97.9 | 112 | 79 | 3 | 40 | 10-124 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM IV
DIESEL RANGE ORGANICS METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: MB 240-206884/23-A

Matrix: Solid Date Extracted: 11/16/2015 08:33

Lab File ID:(1) P5B11708.D Lab File ID:(2)

Date Analyzed:(1) 11/17/2015 22:08 Date Analyzed:(2)

Instrument ID:(1) A2HP5R Instrument ID:(2)

GC Column:(1) ID: GC Column:(2) ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-206884/24-A | 11/17/2015 22:40 | |
| F-VS-UST-P-1 | 240-57930-1 | 11/18/2015 07:32 | |
| F-VS-UST-P-2 | 240-57930-2 | 11/18/2015 08:04 | |
| F-VS-UST-P-3 | 240-57930-3 | 11/18/2015 08:35 | |
| F-VS-UST-P-4 | 240-57930-4 | 11/18/2015 10:10 | |
| F-VB-UST-P-1 | 240-57930-5 | 11/18/2015 10:47 | |
| F-VB-UST-P-2 | 240-57930-6 | 11/18/2015 11:21 | |
| F-VB-UST-P-3 | 240-57930-7 | 11/18/2015 12:32 | |

FORM IV
DIESEL RANGE ORGANICS METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: MB 240-207121/7-A

Matrix: Solid Date Extracted: 11/17/2015 10:02

Lab File ID:(1) P5B11815.D Lab File ID:(2)

Date Analyzed:(1) 11/19/2015 01:29 Date Analyzed:(2)

Instrument ID:(1) A2HP5R Instrument ID:(2)

GC Column:(1) ID: GC Column:(2) ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-207121/8-A | 11/19/2015 02:01 | |
| F-VB-UST-P-4 | 240-57930-8 | 11/19/2015 02:32 | |
| F-VB-UST-P-5 | 240-57930-9 | 11/19/2015 03:04 | |
| F-VB-UST-P-6 | 240-57930-10 | 11/19/2015 03:36 | |
| F-VS-UST-P-DUP | 240-57930-11 | 11/19/2015 04:07 | |
| F-VS-UST-P-DUP MS | 240-57930-11 MS | 11/19/2015 04:38 | |
| F-VS-UST-P-DUP MSD | 240-57930-11 MSD | 11/19/2015 05:10 | |

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Sample No.: STD3 240-205608/10 Date Analyzed: 11/06/2015 17:59

Instrument ID: A2HP5R GC Column: ID: ()

Lab File ID (Standard): P5B10610.D Heated Purge: (Y/N) N

Calibration ID: 31386

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| INITIAL CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| STD3 240-205608/10 ICRT | | 11/06/2015 17:59 | P5B10610.D | 9.83 | | |

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCV 240-207234/7  Date Analyzed: 11/17/2015 21:37

Instrument ID: A2HP5R  GC Column:  ID: ()

Lab File ID (Standard): P5B11707.D  Heated Purge: (Y/N) N

Calibration ID: 31386

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCV 240-207234/7 CCVRT | | 11/17/2015 21:37 | P5B11707.D | 9.83 | | |
| MB 240-206884/23-A | | 11/17/2015 22:08 | P5B11708.D | 9.83 | | |
| LCS 240-206884/24-A | | 11/17/2015 22:40 | P5B11709.D | 9.83 | | |
| CCV 240-207234/18 | | 11/18/2015 03:22 | P5B11718.D | 9.83 | | |
| 240-57930-1 | F-VS-UST-P-1 | 11/18/2015 07:32 | P5B11726.D | 9.83 | | |
| 240-57930-2 | F-VS-UST-P-2 | 11/18/2015 08:04 | P5B11727.D | 9.83 | | |
| 240-57930-3 | F-VS-UST-P-3 | 11/18/2015 08:35 | P5B11728.D | 9.83 | | |
| CCV 240-207234/30 | | 11/18/2015 09:38 | P5B11730.D | 9.83 | | |
| 240-57930-4 | F-VS-UST-P-4 | 11/18/2015 10:10 | P5B11731.D | 9.83 | | |
| 240-57930-5 | F-VB-UST-P-1 | 11/18/2015 10:47 | P5B11732.D | 9.83 | | |
| 240-57930-6 | F-VB-UST-P-2 | 11/18/2015 11:21 | P5B11733.D | 9.83 | | |
| 240-57930-7 | F-VB-UST-P-3 | 11/18/2015 12:32 | P5B11736.D | 9.83 | | |
| CCV 240-207234/38 | | 11/18/2015 13:04 | P5B11737.D | 9.83 | | |

OTPH = o-Terphenyl

OTPH RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Sample No.: CCV 240-207465/7 Date Analyzed: 11/18/2015 21:17

Instrument ID: A2HP5R GC Column: ID: ()

Lab File ID (Standard): P5B11807.D Heated Purge: (Y/N) N

Calibration ID: 31386

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCV 240-207465/7 CCVRT | | 11/18/2015 21:17 | P5B11807.D | 9.83 | | |
| CCV 240-207465/14 | | 11/19/2015 00:58 | P5B11814.D | 9.83 | | |
| MB 240-207121/7-A | | 11/19/2015 01:29 | P5B11815.D | 9.83 | | |
| LCS 240-207121/8-A | | 11/19/2015 02:01 | P5B11816.D | 9.83 | | |
| 240-57930-8 | F-VB-UST-P-4 | 11/19/2015 02:32 | P5B11817.D | 9.83 | | |
| 240-57930-9 | F-VB-UST-P-5 | 11/19/2015 03:04 | P5B11818.D | 9.83 | | |
| 240-57930-10 | F-VB-UST-P-6 | 11/19/2015 03:36 | P5B11819.D | 9.83 | | |
| 240-57930-11 | F-VS-UST-P-DUP | 11/19/2015 04:07 | P5B11820.D | 9.83 | | |
| 240-57930-11 MS | F-VS-UST-P-DUP MS | 11/19/2015 04:38 | P5B11821.D | 9.83 | | |
| 240-57930-11 MSD | F-VS-UST-P-DUP MSD | 11/19/2015 05:10 | P5B11822.D | 9.83 | | |
| CCV 240-207465/23 | | 11/19/2015 05:42 | P5B11823.D | 9.83 | | |

OTPH = o-Terphenyl

OTPH RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-1 | Lab Sample ID: 240-57930-1

Matrix: Solid | Lab File ID: P5B11726.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:30

Extraction Method: 3540C | Date Extracted: 11/16/2015 08:33

Sample wt/vol: 30.28(g) | Date Analyzed: 11/18/2015 07:32

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207234 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 80 | | 19 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 82 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: P5B11727.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:35

Extraction Method: 3540C | Date Extracted: 11/16/2015 08:33

Sample wt/vol: 29.96(g) | Date Analyzed: 11/18/2015 08:04

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207234 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 130 | | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 95 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-3 | Lab Sample ID: 240-57930-3

Matrix: Solid | Lab File ID: P5B11728.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:40

Extraction Method: 3540C | Date Extracted: 11/16/2015 08:33

Sample wt/vol: 30.34(g) | Date Analyzed: 11/18/2015 08:35

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207234 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 42 | | 20 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 75 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-4 | Lab Sample ID: 240-57930-4

Matrix: Solid | Lab File ID: P5B11731.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:45

Extraction Method: 3540C | Date Extracted: 11/16/2015 08:33

Sample wt/vol: 30.19(g) | Date Analyzed: 11/18/2015 10:10

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207234 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 120 | | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 82 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 | Lab Sample ID: 240-57930-5

Matrix: Solid | Lab File ID: P5B11732.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:00

Extraction Method: 3540C | Date Extracted: 11/16/2015 08:33

Sample wt/vol: 29.96(g) | Date Analyzed: 11/18/2015 10:47

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207234 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.6 | U | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 76 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-2 | Lab Sample ID: 240-57930-6

Matrix: Solid | Lab File ID: P5B11733.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:05

Extraction Method: 3540C | Date Extracted: 11/16/2015 08:33

Sample wt/vol: 29.88(g) | Date Analyzed: 11/18/2015 11:21

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207234 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.6 | U | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 79 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-3 | Lab Sample ID: 240-57930-7

Matrix: Solid | Lab File ID: P5B11736.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:10

Extraction Method: 3540C | Date Extracted: 11/16/2015 08:33

Sample wt/vol: 30.03(g) | Date Analyzed: 11/18/2015 12:32

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 19.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207234 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.9 | U | 21 | 7.9 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 69 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-4 | Lab Sample ID: 240-57930-8

Matrix: Solid | Lab File ID: P5B11817.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:15

Extraction Method: 3540C | Date Extracted: 11/17/2015 10:02

Sample wt/vol: 30.20(g) | Date Analyzed: 11/19/2015 02:32

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207465 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.6 | U | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 90 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-5 | Lab Sample ID: 240-57930-9

Matrix: Solid | Lab File ID: P5B11818.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:20

Extraction Method: 3540C | Date Extracted: 11/17/2015 10:02

Sample wt/vol: 29.52(g) | Date Analyzed: 11/19/2015 03:04

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207465 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.8 | U | 20 | 7.8 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 81 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-6 | Lab Sample ID: 240-57930-10

Matrix: Solid | Lab File ID: P5B11819.D

Analysis Method: 8015C | Date Collected: 11/13/2015 10:25

Extraction Method: 3540C | Date Extracted: 11/17/2015 10:02

Sample wt/vol: 30.04(g) | Date Analyzed: 11/19/2015 03:36

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207465 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 8.0 | J | 20 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 83 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP | Lab Sample ID: 240-57930-11

Matrix: Solid | Lab File ID: P5B11820.D

Analysis Method: 8015C | Date Collected: 11/13/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/17/2015 10:02

Sample wt/vol: 30.26(g) | Date Analyzed: 11/19/2015 04:07

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207465 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 51 | | 20 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 79 | | 40-160 |

FORM VI

DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA

RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 205608

SDG No.:

Instrument ID: A2HP5R GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 11/06/2015 15:50 Calibration End Date: 11/06/2015 19:01 Calibration ID: 31386

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-205608/8 | P5B10608.D |
| Level 2 | STD2 240-205608/9 | P5B10609.D |
| Level 3 | STD3 240-205608/10 | P5B10610.D |
| Level 4 | STD4 240-205608/11 | P5B10611.D |
| Level 5 | STD5 240-205608/12 | P5B10612.D |

| ANALYTE | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | | | | | RT WINDOW | AVG RT |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | 9.331 | 9.331 | 9.331 | 9.331 | 9.331 | | | | | | 5.131 - 13.532 | 9.331 |
| o-Terphenyl (Surr) | 9.828 | 9.831 | 9.829 | 9.830 | 9.833 | | | | | | 9.728 - 9.928 | 9.830 |

FORM VI
DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 205608

SDG No.:

Instrument ID: A2HP5R GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 11/06/2015 15:50 Calibration End Date: 11/06/2015 19:01 Calibration ID: 31386

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-205608/8 | P5B10608.D |
| Level 2 | STD2 240-205608/9 | P5B10609.D |
| Level 3 | STD3 240-205608/10 | P5B10610.D |
| Level 4 | STD4 240-205608/11 | P5B10611.D |
| Level 5 | STD5 240-205608/12 | P5B10612.D |

| ANALYTE | CF | | | | CURVE TYPE | COEFFICIENT | | | # | MIN CF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 5 | LVL 2 | LVL 3 | LVL 4 | | B | M1 | M2 | | | | | | | | |
| Diesel Range Organics [C10 - C28] | 156067 174183 | 161462 | 135888 | 139742 | Ave | | 153468.493 | | | | 10.3 | | 20.0 | | | |
| o-Terphenyl (Surr) | 213204 240558 | 209043 | 182141 | 189552 | Ave | | 206899.578 | | | | 11.1 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 205608

SDG No.:

Instrument ID: A2HP5R GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 11/06/2015 15:50 Calibration End Date: 11/06/2015 19:01 Calibration ID: 31386

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-205608/8 | P5B10608.D |
| Level 2 | STD2 240-205608/9 | P5B10609.D |
| Level 3 | STD3 240-205608/10 | P5B10610.D |
| Level 4 | STD4 240-205608/11 | P5B10611.D |
| Level 5 | STD5 240-205608/12 | P5B10612.D |

| ANALYTE | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|
| | | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Diesel Range Organics [C10 - C28] | Ave | 15610573 | 32300497 | 67960720 | 139776482 | 348453260 | 100 | 200 | 500 | 1000 | 2000 |
| o-Terphenyl (Surr) | Ave | 1705635 | 3344683 | 7285634 | 15164183 | 38489230 | 8.00 | 16.0 | 40.0 | 80.0 | 160 |

Curve Type Legend:
Ave = Average

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-205608/13 Calibration Date: 11/06/2015 19:33

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B10613.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 164201 | | 537 | 502 | 7.0 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-205608/13 Calibration Date: 11/06/2015 19:33

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B10613.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207234/7 Calibration Date: 11/17/2015 21:37

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11707.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 150001 | | 489 | 500 | -2.3 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 220429 | | 42.6 | 40.0 | 6.5 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207234/7 Calibration Date: 11/17/2015 21:37

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11707.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207234/18  Calibration Date: 11/18/2015 03:22

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11718.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 163566 | | 533 | 500 | 6.6 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 224743 | | 43.4 | 40.0 | 8.6 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207234/18 Calibration Date: 11/18/2015 03:22

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11718.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207234/30 Calibration Date: 11/18/2015 09:38

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11730.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 168725 | | 550 | 500 | 9.9 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 232043 | | 44.9 | 40.0 | 12.2 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207234/30 Calibration Date: 11/18/2015 09:38

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11730.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207234/38 Calibration Date: 11/18/2015 13:04

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11737.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 164949 | | 538 | 500 | 7.5 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 233758 | | 45.2 | 40.0 | 13.0 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207234/38 Calibration Date: 11/18/2015 13:04

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11737.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207465/7 Calibration Date: 11/18/2015 21:17

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11807.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 179371 | | 585 | 500 | 16.9 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 232427 | | 44.9 | 40.0 | 12.3 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207465/7 Calibration Date: 11/18/2015 21:17

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11807.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207465/14 Calibration Date: 11/19/2015 00:58

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11814.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 174581 | | 569 | 500 | 13.8 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 215106 | | 41.6 | 40.0 | 4.0 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207465/14 Calibration Date: 11/19/2015 00:58

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11814.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207465/23 Calibration Date: 11/19/2015 05:42

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11823.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 171249 | | 558 | 500 | 11.6 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 229002 | | 44.3 | 40.0 | 10.7 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207465/23 Calibration Date: 11/19/2015 05:42

Instrument ID: A2HP5R Calib Start Date: 11/06/2015 15:50

GC Column: ID: Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11823.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.13 | 13.53 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206884/23-A

Matrix: Solid | Lab File ID: P5B11708.D

Analysis Method: 8015C | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/16/2015 08:33

Sample wt/vol: 30(g) | Date Analyzed: 11/17/2015 22:08

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207234 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 6.4 | U | 17 | 6.4 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 77 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-207121/7-A

Matrix: Solid | Lab File ID: P5B11815.D

Analysis Method: 8015C | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/17/2015 10:02

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 01:29

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207465 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 6.4 | U | 17 | 6.4 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 75 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206884/24-A

Matrix: Solid | Lab File ID: P5B11709.D

Analysis Method: 8015C | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/16/2015 08:33

Sample wt/vol: 30(g) | Date Analyzed: 11/17/2015 22:40

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207234 | Units: mg/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 82 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-207121/8-A

Matrix: Solid | Lab File ID: P5B11816.D

Analysis Method: 8015C | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/17/2015 10:02

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 02:01

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207465 | Units: mg/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 103 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP MS | Lab Sample ID: 240-57930-11 MS

Matrix: Solid | Lab File ID: P5B11821.D

Analysis Method: 8015C | Date Collected: 11/13/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/17/2015 10:02

Sample wt/vol: 29.91(g) | Date Analyzed: 11/19/2015 04:38

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207465 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 124 | | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 93 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP MSD | Lab Sample ID: 240-57930-11 MSD

Matrix: Solid | Lab File ID: P5B11822.D

Analysis Method: 8015C | Date Collected: 11/13/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/17/2015 10:02

Sample wt/vol: 30.50(g) | Date Analyzed: 11/19/2015 05:10

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207465 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 126 | | 19 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 93 | | 40-160 |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP5R Start Date: 11/06/2015 15:50

Analysis Batch Number: 205608 End Date: 11/06/2015 19:33

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-205608/8 IC | | 11/06/2015 15:50 | 1 | P5B10608.D | |
| STD2 240-205608/9 IC | | 11/06/2015 17:28 | 1 | P5B10609.D | |
| STD3 240-205608/10 ICRT | | 11/06/2015 17:59 | 1 | P5B10610.D | |
| STD4 240-205608/11 IC | | 11/06/2015 18:30 | 1 | P5B10611.D | |
| STD5 240-205608/12 IC | | 11/06/2015 19:01 | 1 | P5B10612.D | |
| ICV 240-205608/13 | | 11/06/2015 19:33 | 1 | P5B10613.D | |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP5R Start Date: 11/17/2015 21:37

Analysis Batch Number: 207234 End Date: 11/18/2015 13:04

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-207234/7 CCVRT | | 11/17/2015 21:37 | 1 | P5B11707.D | |
| MB 240-206884/23-A | | 11/17/2015 22:08 | 1 | P5B11708.D | |
| LCS 240-206884/24-A | | 11/17/2015 22:40 | 1 | P5B11709.D | |
| CCV 240-207234/18 | | 11/18/2015 03:22 | 1 | P5B11718.D | |
| 240-57930-1 | F-VS-UST-P-1 | 11/18/2015 07:32 | 1 | P5B11726.D | |
| 240-57930-2 | F-VS-UST-P-2 | 11/18/2015 08:04 | 1 | P5B11727.D | |
| 240-57930-3 | F-VS-UST-P-3 | 11/18/2015 08:35 | 1 | P5B11728.D | |
| CCV 240-207234/30 | | 11/18/2015 09:38 | 1 | P5B11730.D | |
| 240-57930-4 | F-VS-UST-P-4 | 11/18/2015 10:10 | 1 | P5B11731.D | |
| 240-57930-5 | F-VB-UST-P-1 | 11/18/2015 10:47 | 1 | P5B11732.D | |
| 240-57930-6 | F-VB-UST-P-2 | 11/18/2015 11:21 | 1 | P5B11733.D | |
| 240-57930-7 | F-VB-UST-P-3 | 11/18/2015 12:32 | 1 | P5B11736.D | |
| CCV 240-207234/38 | | 11/18/2015 13:04 | 1 | P5B11737.D | |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP5R Start Date: 11/18/2015 21:17

Analysis Batch Number: 207465 End Date: 11/19/2015 05:42

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-207465/7 CCVRT | | 11/18/2015 21:17 | 1 | P5B11807.D | |
| ZZZZZ | | 11/18/2015 21:49 | 1 | | |
| ZZZZZ | | 11/18/2015 22:20 | 1 | | |
| ZZZZZ | | 11/18/2015 22:52 | 1 | | |
| ZZZZZ | | 11/18/2015 23:23 | 1 | | |
| ZZZZZ | | 11/18/2015 23:55 | 1 | | |
| ZZZZZ | | 11/19/2015 00:26 | 1 | | |
| CCV 240-207465/14 | | 11/19/2015 00:58 | 1 | P5B11814.D | |
| MB 240-207121/7-A | | 11/19/2015 01:29 | 1 | P5B11815.D | |
| LCS 240-207121/8-A | | 11/19/2015 02:01 | 1 | P5B11816.D | |
| 240-57930-8 | F-VB-UST-P-4 | 11/19/2015 02:32 | 1 | P5B11817.D | |
| 240-57930-9 | F-VB-UST-P-5 | 11/19/2015 03:04 | 1 | P5B11818.D | |
| 240-57930-10 | F-VB-UST-P-6 | 11/19/2015 03:36 | 1 | P5B11819.D | |
| 240-57930-11 | F-VS-UST-P-DUP | 11/19/2015 04:07 | 1 | P5B11820.D | |
| 240-57930-11 MS | F-VS-UST-P-DUP MS | 11/19/2015 04:38 | 1 | P5B11821.D | |
| 240-57930-11 MSD | F-VS-UST-P-DUP MSD | 11/19/2015 05:10 | 1 | P5B11822.D | |
| CCV 240-207465/23 | | 11/19/2015 05:42 | 1 | P5B11823.D | |

# DIESEL RANGE ORGANICS BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 206884  Batch Start Date: 11/16/15 11:50  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/17/15 03:50

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exDIESELspkw 00068 | exOTPSURRW 00006 | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57930-D-1 | F-VS-UST-P-1 | 3540C, 8015C | T | 30.28 g | 5 mL | | 1 mL | | |
| 240-57930-D-2 | F-VS-UST-P-2 | 3540C, 8015C | T | 29.96 g | 5 mL | | 1 mL | | |
| 240-57930-D-3 | F-VS-UST-P-3 | 3540C, 8015C | T | 30.34 g | 5 mL | | 1 mL | | |
| 240-57930-D-4 | F-VS-UST-P-4 | 3540C, 8015C | T | 30.19 g | 5 mL | | 1 mL | | |
| 240-57930-D-5 | F-VB-UST-P-1 | 3540C, 8015C | T | 29.96 g | 5 mL | | 1 mL | | |
| 240-57930-D-6 | F-VB-UST-P-2 | 3540C, 8015C | T | 29.88 g | 5 mL | | 1 mL | | |
| 240-57930-D-7 | F-VB-UST-P-3 | 3540C, 8015C | T | 30.03 g | 5 mL | | 1 mL | | |
| MB 240-206884/23 | | 3540C, 8015C | | 30 g | 5 mL | | 1 mL | | |
| LCS 240-206884/24 | | 3540C, 8015C | | 30 g | 5 mL | 1 mL | 1 mL | | |

| Batch Notes | |
|---|---|
| Balance ID | 1339907 |
| Na2SO4 Lot Number | 2204509 |
| Prep Solvent Lot # | 2287347 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# DIESEL RANGE ORGANICS BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 207121  Batch Start Date: 11/17/15 12:05  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/18/15 04:05

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exDIESELspkw 00068 | exOTPSURRW 00006 | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57930-D-8 | F-VB-UST-P-4 | 3540C, 8015C | T | 30.20 g | 5 mL | | 1 mL | | |
| 240-57930-D-9 | F-VB-UST-P-5 | 3540C, 8015C | T | 29.52 g | 5 mL | | 1 mL | | |
| 240-57930-D-10 | F-VB-UST-P-6 | 3540C, 8015C | T | 30.04 g | 5 mL | | 1 mL | | |
| 240-57930-D-11 | F-VS-UST-P-DUP | 3540C, 8015C | T | 30.26 g | 5 mL | | 1 mL | | |
| 240-57930-D-11 MS | F-VS-UST-P-DUP | 3540C, 8015C | T | 29.91 g | 5 mL | 1 mL | 1 mL | | |
| 240-57930-D-11 MSD | F-VS-UST-P-DUP | 3540C, 8015C | T | 30.50 g | 5 mL | 1 mL | 1 mL | | |
| MB 240-207121/7 | | 3540C, 8015C | | 30 g | 5 mL | | 1 mL | | |
| LCS 240-207121/8 | | 3540C, 8015C | | 30 g | 5 mL | 1 mL | 1 mL | | |

| Batch Notes | |
|---|---|
| Balance ID | 1339912 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2287348 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8082A

## Polychlorinated Biphenyls (PCBs) (GC) by Method 8082A

FORM II
PCBS SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low

GC Column (2): CLP-1 ID: 0.53(mm)

| Client Sample ID | Lab Sample ID | TCX2 # | DCB2 # |
|---|---|---|---|
| F-VS-UST-P-1 | 240-57930-1 | 83 | 72 |
| F-VS-UST-P-2 | 240-57930-2 | 61 | 70 |
| F-VS-UST-P-3 | 240-57930-3 | 67 | 90 |
| F-VS-UST-P-4 | 240-57930-4 | 78 | 141 |
| F-VB-UST-P-1 | 240-57930-5 | 83 | 82 |
| F-VB-UST-P-2 | 240-57930-6 | 70 | 66 |
| F-VB-UST-P-3 | 240-57930-7 | 67 | 66 |
| F-VB-UST-P-4 | 240-57930-8 | 83 | 79 |
| F-VB-UST-P-5 | 240-57930-9 | 80 | 71 |
| F-VB-UST-P-6 | 240-57930-10 | 72 | 71 |
| F-VS-UST-P-DUP | 240-57930-11 | 82 | 78 |
| | MB 240-206917/21-A | 74 | 79 |
| | LCS 240-206917/22-A | 75 | 77 |
| | 240-57880-E-1-B MS | 75 | 94 |
| | 240-57880-E-1-C MSD | 90 | 87 |

| | QC LIMITS |
|---|---|
| TCX = Tetrachloro-m-xylene | 13-134 |
| DCB = DCB Decachlorobiphenyl | 10-155 |

# Column to be used to flag recovery values

FORM II 8082A

FORM III
PCBS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P1200063.D

Lab ID: LCS 240-206917/22-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Aroclor-1016 | 333 | 251 | 75 | 51-120 | |
| Aroclor-1260 | 334 | 253 | 76 | 48-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM III
PCBS MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P1200066.D

Lab ID: 240-57880-E-1-B MS Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Aroclor-1016 | 443 | 16 U | 518 | 117 | 17-141 | |
| Aroclor-1260 | 444 | 71 | 266 | 44 | 19-137 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM III
PCBS MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: P1200067.D

Lab ID: 240-57880-E-1-C MSD  Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Aroclor-1016 | 446 | 638 | 143 | 18 | 56 | 17-141 | F1 |
| Aroclor-1260 | 447 | 331 | 55 | 21 | 55 | 19-137 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM IV
PCBS METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: MB 240-206917/21-A

Matrix: Solid  Date Extracted: 11/16/2015 09:32

Lab File ID:(1)  Lab File ID:(2) P1200062.D

Date Analyzed:(1)  Date Analyzed:(2) 11/19/2015 01:03

Instrument ID:(1)  Instrument ID:(2) A2HP12

GC Column:(1)  ID:  GC Column:(2) CLP-1  ID: 0.53(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| F-VB-UST-P-2 | 240-57930-6 | | 11/18/2015 22:02 |
| F-VB-UST-P-3 | 240-57930-7 | | 11/18/2015 22:20 |
| F-VB-UST-P-4 | 240-57930-8 | | 11/18/2015 22:38 |
| F-VB-UST-P-5 | 240-57930-9 | | 11/18/2015 22:56 |
| F-VB-UST-P-6 | 240-57930-10 | | 11/18/2015 23:15 |
| F-VS-UST-P-DUP | 240-57930-11 | | 11/18/2015 23:33 |
| F-VS-UST-P-1 | 240-57930-1 | | 11/18/2015 23:51 |
| F-VS-UST-P-2 | 240-57930-2 | | 11/19/2015 00:09 |
| F-VS-UST-P-3 | 240-57930-3 | | 11/19/2015 00:27 |
| F-VS-UST-P-4 | 240-57930-4 | | 11/19/2015 00:45 |
| | LCS 240-206917/22-A | | 11/19/2015 01:21 |
| F-VB-UST-P-1 | 240-57930-5 | | 11/19/2015 01:58 |
| | 240-57880-E-1-B MS | | 11/19/2015 02:16 |
| | 240-57880-E-1-C MSD | | 11/19/2015 02:34 |

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: STD05 240-207007/31  Date Analyzed: 11/17/2015 04:23

Instrument ID: A2HP12  GC Column: CLP-1  ID: 0.53(mm)

Lab File ID (Standard): P1200066.D  Heated Purge: (Y/N) N

Calibration ID: 31555

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 188112954 | 2.68 | | | | |
| UPPER LIMIT | | 376225908 | 3.18 | | | | |
| LOWER LIMIT | | 94056477 | 2.18 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-207007/39 | | 184850376 | 2.69 | | | | |
| CCV 240-207364/3 CCVIS | | 190140766 | 2.68 | | | | |

BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Sample No.: CCV 240-207364/3 Date Analyzed: 11/18/2015 18:06

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm)

Lab File ID (Standard): P1200039.D Heated Purge: (Y/N) N

Calibration ID: 31555

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 190140766 | 2.68 | | | | |
| UPPER LIMIT | | 380281532 | 3.18 | | | | |
| LOWER LIMIT | | 95070383 | 2.18 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-207364/4 | | 188081408 | 2.68 | | | | |
| CCV 240-207364/5 | | 202072165 | 2.68 | | | | |
| CCV 240-207364/6 | | 192999673 | 2.69 | | | | |
| CCV 240-207364/7 | | 187722278 | 2.68 | | | | |
| 240-57930-6 | F-VB-UST-P-2 | 206915927 | 2.69 | | | | |
| 240-57930-7 | F-VB-UST-P-3 | 206112342 | 2.69 | | | | |
| 240-57930-8 | F-VB-UST-P-4 | 204369410 | 2.69 | | | | |
| 240-57930-9 | F-VB-UST-P-5 | 215387279 | 2.68 | | | | |
| 240-57930-10 | F-VB-UST-P-6 | 216471663 | 2.68 | | | | |
| 240-57930-11 | F-VS-UST-P-DUP | 216144605 | 2.69 | | | | |
| 240-57930-1 | F-VS-UST-P-1 | 208154897 | 2.69 | | | | |
| 240-57930-2 | F-VS-UST-P-2 | 205161190 | 2.69 | | | | |
| 240-57930-3 | F-VS-UST-P-3 | 213739691 | 2.68 | | | | |
| 240-57930-4 | F-VS-UST-P-4 | 218600878 | 2.68 | | | | |
| MB 240-206917/21-A | | 207945676 | 2.68 | | | | |
| LCS 240-206917/22-A | | 205793008 | 2.69 | | | | |

BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Sample No.: CCVIS 240-207364/28  Date Analyzed: 11/19/2015 01:40

Instrument ID: A2HP12  GC Column: CLP-1  ID: 0.53(mm)

Lab File ID (Standard): P1200064.D  Heated Purge: (Y/N) N

Calibration ID: 31555

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 203830640 | 2.68 | | | | |
| UPPER LIMIT | | 407661280 | 3.18 | | | | |
| LOWER LIMIT | | 101915320 | 2.18 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| 240-57930-5 | F-VB-UST-P-1 | 210716095 | 2.69 | | | | |
| 240-57880-E-1-B MS | | 219185526 | 2.68 | | | | |
| 240-57880-E-1-C MSD | | 210947173 | 2.68 | | | | |

BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM X
IDENTIFICATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Client Sample ID: Lab Sample ID: 240-57880-E-1-B MS

Instrument ID (1): Instrument ID (2): A2HP12

Date Analyzed (1): Date Analyzed (2): 11/19/2015 02:16

GC Column (1): ID: GC Column (2): CLP-1 ID: 0.53(mm)

| ANALYTE | COL | PEAK | RT | RT WINDOW | | CONCENTRATION | | |
|---|---|---|---|---|---|---|---|---|
| | | | | FROM | TO | PEAK | MEAN | |
| Aroclor-1016 | 1 | 1 | 4.61 | 4.58 | 4.64 | 581 | 518 | |
| | | 2 | 5.11 | 5.08 | 5.14 | 594 | | |
| | | 3 | 5.74 | 5.71 | 5.77 | 517 | | |
| | | 4 | 5.91 | 5.88 | 5.94 | 419 | | |
| | | 5 | 6.45 | 6.42 | 6.48 | 482 | | |
| Aroclor-1260 | 1 | 1 | 7.95 | 7.93 | 7.99 | 275 | 266 | |
| | | 2 | 8.29 | 8.26 | 8.32 | 240 | | |
| | | 3 | 8.61 | 8.58 | 8.64 | 249 | | |
| | | 4 | 9.40 | 9.37 | 9.43 | 281 | | |
| | | 5 | 9.72 | 9.69 | 9.75 | 284 | | |

FORM X
IDENTIFICATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Client Sample ID: Lab Sample ID: 240-57880-E-1-C MSD

Instrument ID (1): Instrument ID (2): A2HP12

Date Analyzed (1): Date Analyzed (2): 11/19/2015 02:34

GC Column (1): ID: GC Column (2): CLP-1 ID: 0.53(mm)

| ANALYTE | COL | PEAK | RT | RT WINDOW | | CONCENTRATION | | |
|---|---|---|---|---|---|---|---|---|
| | | | | FROM | TO | PEAK | MEAN | |
| Aroclor-1016 | 1 | 1 | 4.61 | 4.58 | 4.64 | 693 | 638 | |
| | | 2 | 5.11 | 5.08 | 5.14 | 722 | | |
| | | 3 | 5.74 | 5.71 | 5.77 | 636 | | |
| | | 4 | 5.91 | 5.88 | 5.94 | 544 | | |
| | | 5 | 6.45 | 6.42 | 6.48 | 597 | | |
| Aroclor-1260 | 1 | 1 | 7.95 | 7.93 | 7.99 | 341 | 331 | |
| | | 2 | 8.29 | 8.26 | 8.32 | 298 | | |
| | | 3 | 8.61 | 8.58 | 8.64 | 297 | | |
| | | 4 | 9.40 | 9.37 | 9.43 | 363 | | |
| | | 5 | 9.72 | 9.69 | 9.75 | 356 | | |

FORM X
IDENTIFICATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Client Sample ID: Lab Sample ID: LCS 240-206917/22-A

Instrument ID (1): Instrument ID (2): A2HP12

Date Analyzed (1): Date Analyzed (2): 11/19/2015 01:21

GC Column (1): ID: GC Column (2): CLP-1 ID: 0.53(mm)

| ANALYTE | COL | PEAK | RT | RT WINDOW | | CONCENTRATION | | |
|---|---|---|---|---|---|---|---|---|
| | | | | FROM | TO | PEAK | MEAN | |
| Aroclor-1016 | 1 | 1 | 4.61 | 4.58 | 4.64 | 244 | 251 | |
| | | 2 | 5.11 | 5.08 | 5.14 | 248 | | |
| | | 3 | 5.74 | 5.71 | 5.77 | 257 | | |
| | | 4 | 5.92 | 5.88 | 5.94 | 256 | | |
| | | 5 | 6.45 | 6.42 | 6.48 | 250 | | |
| Aroclor-1260 | 1 | 1 | 7.96 | 7.92 | 7.98 | 260 | 253 | |
| | | 2 | 8.29 | 8.26 | 8.32 | 259 | | |
| | | 3 | 8.61 | 8.58 | 8.64 | 250 | | |
| | | 4 | 9.40 | 9.37 | 9.43 | 258 | | |
| | | 5 | 9.72 | 9.69 | 9.75 | 240 | | |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-1 | Lab Sample ID: 240-57930-1

Matrix: Solid | Lab File ID: P1200058.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:30

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 29.62(g) | Date Analyzed: 11/18/2015 23:51

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 83 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 72 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-2 | Lab Sample ID: 240-57930-2

Matrix: Solid | Lab File ID: P1200059.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:35

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 30.44(g) | Date Analyzed: 11/19/2015 00:09

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 61 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 70 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-3 | Lab Sample ID: 240-57930-3

Matrix: Solid | Lab File ID: P1200060.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:40

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 29.65(g) | Date Analyzed: 11/19/2015 00:27

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.6 | U | 39 | 9.6 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 67 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 90 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-4 | Lab Sample ID: 240-57930-4

Matrix: Solid | Lab File ID: P1200061.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:45

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 30.41(g) | Date Analyzed: 11/19/2015 00:45

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 16.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 78 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 141 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-1 | Lab Sample ID: 240-57930-5

Matrix: Solid | Lab File ID: P1200065.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:00

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 30.45(g) | Date Analyzed: 11/19/2015 01:58

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 15.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 83 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 82 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-2 | Lab Sample ID: 240-57930-6

Matrix: Solid | Lab File ID: P1200052.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:05

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 29.88(g) | Date Analyzed: 11/18/2015 22:02

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 15.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 70 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 66 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-3 | Lab Sample ID: 240-57930-7

Matrix: Solid | Lab File ID: P1200053.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:10

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 30.01(g) | Date Analyzed: 11/18/2015 22:20

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 19.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 15 | U | 41 | 15 |
| 11104-28-2 | Aroclor-1221 | 20 | U | 41 | 20 |
| 11141-16-5 | Aroclor-1232 | 25 | U | 41 | 25 |
| 53469-21-9 | Aroclor-1242 | 14 | U | 41 | 14 |
| 12672-29-6 | Aroclor-1248 | 9.9 | U | 41 | 9.9 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 41 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 41 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 67 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 66 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-4 | Lab Sample ID: 240-57930-8

Matrix: Solid | Lab File ID: P1200054.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:15

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 29.74(g) | Date Analyzed: 11/18/2015 22:38

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 40 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.6 | U | 40 | 9.6 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 83 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 79 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-5 | Lab Sample ID: 240-57930-9

Matrix: Solid | Lab File ID: P1200055.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:20

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 29.69(g) | Date Analyzed: 11/18/2015 22:56

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 16.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 15 | U | 40 | 15 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.7 | U | 40 | 9.7 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 80 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 71 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VB-UST-P-6 | Lab Sample ID: 240-57930-10

Matrix: Solid | Lab File ID: P1200056.D

Analysis Method: 8082A | Date Collected: 11/13/2015 10:25

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 29.68(g) | Date Analyzed: 11/18/2015 23:15

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 72 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 71 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: F-VS-UST-P-DUP | Lab Sample ID: 240-57930-11

Matrix: Solid | Lab File ID: P1200057.D

Analysis Method: 8082A | Date Collected: 11/13/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 29.83(g) | Date Analyzed: 11/18/2015 23:33

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 82 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 78 | | 10-155 |

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 20:14 Calibration End Date: 11/16/2015 21:45 Calibration ID: 31527

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/4 | P1200039.D |
| Level 2 | STD01 240-207007/5 | P1200040.D |
| Level 3 | STD02 240-207007/6 | P1200041.D |
| Level 4 | STD05 240-207007/7 | P1200042.D |
| Level 5 | STD1 240-207007/8 | P1200043.D |
| Level 6 | STD2 240-207007/9 | P1200044.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1232 Peak 1 | 0.0146 0.0120 | 0.0146 | 0.0137 | 0.0137 | 0.0123 | Lin1 | 0.0002 | 0.0122 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 2 | 0.0104 0.0088 | 0.0106 | 0.0101 | 0.0100 | 0.0090 | Lin1 | 0.0001 | 0.0089 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 3 | 0.0214 0.0206 | 0.0218 | 0.0204 | 0.0219 | 0.0205 | Lin1 | 0.0001 | 0.0207 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1232 Peak 4 | 0.0088 0.0084 | 0.0090 | 0.0081 | 0.0093 | 0.0085 | Lin1 | 0 | 0.0085 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 5 | 0.0088 0.0076 | 0.0092 | 0.0087 | 0.0086 | 0.0078 | Lin1 | 0.0001 | 0.0078 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1262 Peak 1 | 0.0288 0.0238 | 0.0249 | 0.0236 | 0.0241 | 0.0228 | Lin1 | 0.0002 | 0.0233 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1262 Peak 2 | 0.0420 0.0453 | 0.0443 | 0.0426 | 0.0462 | 0.0441 | Lin1 | 0 | 0.0452 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 3 | 0.0380 0.0415 | 0.0404 | 0.0397 | 0.0426 | 0.0405 | Lin1 | 0 | 0.0415 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 4 | 0.0853 0.0976 | 0.0928 | 0.0961 | 0.0994 | 0.0962 | Lin1 | -0.001 | 0.0979 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 5 | 0.0364 0.0385 | 0.0386 | 0.0374 | 0.0395 | 0.0372 | Lin1 | 0 | 0.0383 | | | | | | | 0.9990 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 20:14 Calibration End Date: 11/16/2015 21:45 Calibration ID: 31527

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/4 | P1200039.D |
| Level 2 | STD01 240-207007/5 | P1200040.D |
| Level 3 | STD02 240-207007/6 | P1200041.D |
| Level 4 | STD05 240-207007/7 | P1200042.D |
| Level 5 | STD1 240-207007/8 | P1200043.D |
| Level 6 | STD2 240-207007/9 | P1200044.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1232 Peak 1 | BNB | Lin1 | 2864462<br>94613461 | 5743911 | 10579052 | 25148552 | 48682811 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 2 | BNB | Lin1 | 2042318<br>69008204 | 4176797 | 7799621 | 18338381 | 35520928 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 3 | BNB | Lin1 | 4204092<br>162234139 | 8567857 | 15765426 | 40096819 | 80981509 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 4 | BNB | Lin1 | 1732949<br>65818756 | 3551392 | 6289163 | 17045553 | 33804666 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 5 | BNB | Lin1 | 1725581<br>59962732 | 3622594 | 6723266 | 15858414 | 31021460 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 1 | BNB | Lin1 | 5662298<br>186973877 | 9773527 | 18252707 | 44264730 | 89983608 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 2 | BNB | Lin1 | 8249286<br>356026718 | 17431705 | 32928458 | 84853483 | 174547806 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 3 | BNB | Lin1 | 7469803<br>326504754 | 15875631 | 30684939 | 78118840 | 160322518 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 4 | BNB | Lin1 | 16781995<br>766924713 | 36487355 | 74250353 | 182460886 | 380317708 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 5 | BNB | Lin1 | 7158889<br>302670598 | 15182112 | 28944443 | 72530686 | 147149653 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 22:03 Calibration End Date: 11/16/2015 23:33 Calibration ID: 31533

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/10 | P1200045.D |
| Level 2 | STD01 240-207007/11 | P1200046.D |
| Level 3 | STD02 240-207007/12 | P1200047.D |
| Level 4 | STD05 240-207007/13 | P1200048.D |
| Level 5 | STD1 240-207007/14 | P1200049.D |
| Level 6 | STD2 240-207007/15 | P1200050.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1242 Peak 1 | 0.0115 0.0082 | 0.0100 | 0.0094 | 0.0091 | 0.0087 | Lin1 | 0.0002 | 0.0083 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 2 | 0.0201 0.0153 | 0.0180 | 0.0172 | 0.0169 | 0.0161 | Lin1 | 0.0003 | 0.0155 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 3 | 0.0408 0.0378 | 0.0334 | 0.0373 | 0.0380 | 0.0380 | Lin1 | 0 | 0.0379 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 4 | 0.0176 0.0154 | 0.0131 | 0.0162 | 0.0160 | 0.0157 | Lin1 | 0 | 0.0155 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1242 Peak 5 | 0.0196 0.0154 | 0.0173 | 0.0163 | 0.0163 | 0.0157 | Lin1 | 0.0002 | 0.0155 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 1 | 0.1160 0.1140 | 0.1051 | 0.1050 | 0.1121 | 0.1102 | Lin1 | 0 | 0.1126 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1268 Peak 2 | 0.1129 0.1102 | 0.1035 | 0.1031 | 0.1082 | 0.1086 | Lin1 | 0 | 0.1094 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 3 | 0.0941 0.0941 | 0.0865 | 0.0873 | 0.0935 | 0.0918 | Lin1 | 0 | 0.0934 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 4 | 0.0427 0.0386 | 0.0375 | 0.0368 | 0.0384 | 0.0379 | Lin1 | 0.0001 | 0.0382 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 5 | 0.3010 0.3253 | 0.2825 | 0.2850 | 0.3104 | 0.3105 | Lin1 | -0.003 | 0.3204 | | | | | | | 0.9990 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 22:03 Calibration End Date: 11/16/2015 23:33 Calibration ID: 31533

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/10 | P1200045.D |
| Level 2 | STD01 240-207007/11 | P1200046.D |
| Level 3 | STD02 240-207007/12 | P1200047.D |
| Level 4 | STD05 240-207007/13 | P1200048.D |
| Level 5 | STD1 240-207007/14 | P1200049.D |
| Level 6 | STD2 240-207007/15 | P1200050.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1242 Peak 1 | BNB | Lin1 | 2149767<br>64172683 | 3972268 | 7542438 | 18026385 | 33712614 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 2 | BNB | Lin1 | 3777832<br>120095985 | 7167048 | 13831308 | 33547847 | 62739379 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 3 | BNB | Lin1 | 7649911<br>296721464 | 13315192 | 29900733 | 75211737 | 147999390 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 4 | BNB | Lin1 | 3299090<br>120600215 | 5223145 | 12995644 | 31796203 | 60982553 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 5 | BNB | Lin1 | 3672196<br>121118126 | 6903841 | 13110974 | 32386776 | 61129587 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 1 | BNB | Lin1 | 21759307<br>894333238 | 41863019 | 84230109 | 222130676 | 428635453 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 2 | BNB | Lin1 | 21178393<br>864175846 | 41209432 | 82708339 | 214470756 | 422311987 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 3 | BNB | Lin1 | 17646266<br>738380844 | 34461725 | 70030810 | 185306823 | 357127789 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 4 | BNB | Lin1 | 8003779<br>302725873 | 14951207 | 29550936 | 76139152 | 147585310 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 5 | BNB | Lin1 | 56470647<br>2551765384 | 112539547 | 228706125 | 615081152 | 1208047688 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12  GC Column: CLP-1  ID: 0.53(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 23:51  Calibration End Date: 11/17/2015 01:22  Calibration ID: 31539

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/16 | P1200051.D |
| Level 2 | STD01 240-207007/17 | P1200052.D |
| Level 3 | STD02 240-207007/18 | P1200053.D |
| Level 4 | STD05 240-207007/19 | P1200054.D |
| Level 5 | STD1 240-207007/20 | P1200055.D |
| Level 6 | STD2 240-207007/21 | P1200056.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1248 Peak 1 | 0.0111 0.0082 | 0.0099 | 0.0099 | 0.0090 | 0.0084 | Lin1 | 0.0002 | 0.0082 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1248 Peak 2 | 0.0277 0.0242 | 0.0256 | 0.0251 | 0.0243 | 0.0239 | Lin1 | 0.0002 | 0.0240 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 3 | 0.0240 0.0225 | 0.0232 | 0.0233 | 0.0226 | 0.0219 | Lin1 | 0.0001 | 0.0223 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 4 | 0.0199 0.0193 | 0.0194 | 0.0192 | 0.0192 | 0.0190 | Lin1 | 0 | 0.0192 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 5 | 0.0138 0.0128 | 0.0122 | 0.0128 | 0.0125 | 0.0116 | Lin1 | 0 | 0.0124 | | | | | | | 0.9980 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 23:51 Calibration End Date: 11/17/2015 01:22 Calibration ID: 31539

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/16 | P1200051.D |
| Level 2 | STD01 240-207007/17 | P1200052.D |
| Level 3 | STD02 240-207007/18 | P1200053.D |
| Level 4 | STD05 240-207007/19 | P1200054.D |
| Level 5 | STD1 240-207007/20 | P1200055.D |
| Level 6 | STD2 240-207007/21 | P1200056.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1248 Peak 1 | BNB | Lin1 | 2214671<br>65344382 | 3893329 | 7710225 | 17465613 | 33070601 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 2 | BNB | Lin1 | 5540374<br>193008938 | 10066768 | 19570159 | 46979815 | 94056087 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 3 | BNB | Lin1 | 4803789<br>179213673 | 9149673 | 18179130 | 43818711 | 86223823 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 4 | BNB | Lin1 | 3975067<br>153520953 | 7642679 | 14994793 | 37207246 | 74584605 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 5 | BNB | Lin1 | 2764694<br>102029173 | 4797959 | 10009264 | 24248210 | 45732697 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/17/2015 01:40 Calibration End Date: 11/17/2015 03:10 Calibration ID: 31549

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/22 | P1200057.D |
| Level 2 | STD01 240-207007/23 | P1200058.D |
| Level 3 | STD02 240-207007/24 | P1200059.D |
| Level 4 | STD05 240-207007/25 | P1200060.D |
| Level 5 | STD1 240-207007/26 | P1200061.D |
| Level 6 | STD2 240-207007/27 | P1200062.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1221 Peak 1 | 0.0083 0.0057 | 0.0074 | 0.0067 | 0.0065 | 0.0061 | Lin1 | 0.0002 | 0.0058 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1221 Peak 2 | 0.0055 0.0037 | 0.0049 | 0.0045 | 0.0042 | 0.0039 | Lin1 | 0.0001 | 0.0037 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1221 Peak 3 | 0.0208 0.0145 | 0.0183 | 0.0167 | 0.0164 | 0.0154 | Lin1 | 0.0004 | 0.0147 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1254 Peak 1 | 0.0259 0.0202 | 0.0228 | 0.0210 | 0.0212 | 0.0203 | Lin1 | 0.0003 | 0.0201 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 2 | 0.0354 0.0290 | 0.0313 | 0.0287 | 0.0298 | 0.0292 | Lin1 | 0.0003 | 0.0289 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 3 | 0.0466 0.0419 | 0.0442 | 0.0413 | 0.0427 | 0.0391 | Lin1 | 0.0003 | 0.0410 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1254 Peak 4 | 0.0442 0.0360 | 0.0392 | 0.0345 | 0.0358 | 0.0355 | Lin1 | 0.0003 | 0.0354 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1254 Peak 5 | 0.0544 0.0459 | 0.0486 | 0.0448 | 0.0465 | 0.0453 | Lin1 | 0.0003 | 0.0454 | | | | | | | 1.0000 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/17/2015 01:40 Calibration End Date: 11/17/2015 03:10 Calibration ID: 31549

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/22 | P1200057.D |
| Level 2 | STD01 240-207007/23 | P1200058.D |
| Level 3 | STD02 240-207007/24 | P1200059.D |
| Level 4 | STD05 240-207007/25 | P1200060.D |
| Level 5 | STD1 240-207007/26 | P1200061.D |
| Level 6 | STD2 240-207007/27 | P1200062.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1221 Peak 1 | BNB | Lin1 | 1617753<br>43159048 | 2834146 | 5288305 | 12630620 | 23444359 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1221 Peak 2 | BNB | Lin1 | 1073929<br>27877416 | 1853226 | 3515548 | 8239497 | 15092745 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1221 Peak 3 | BNB | Lin1 | 4061658<br>109234720 | 6958986 | 13214576 | 31899214 | 59455965 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 1 | BNB | Lin1 | 5070982<br>152082598 | 8694593 | 16601983 | 41262994 | 78538732 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 2 | BNB | Lin1 | 6915782<br>218988362 | 11916252 | 22685227 | 57976400 | 113180990 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 3 | BNB | Lin1 | 9122282<br>316127715 | 16841606 | 32606706 | 82955431 | 151535520 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 4 | BNB | Lin1 | 8646084<br>271574308 | 14939001 | 27264214 | 69542078 | 137323388 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 5 | BNB | Lin1 | 10643906<br>346633557 | 18530884 | 35338412 | 90448562 | 175425464 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57930-1  Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12  GC Column: CLP-1  ID: 0.53(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/17/2015 03:29  Calibration End Date: 11/17/2015 04:59  Calibration ID: 31555

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/28 | P1200063.D |
| Level 2 | STD01 240-207007/29 | P1200064.D |
| Level 3 | STD02 240-207007/30 | P1200065.D |
| Level 4 | STD05 240-207007/31 | P1200066.D |
| Level 5 | STD1 240-207007/32 | P1200067.D |
| Level 6 | STD2 240-207007/33 | P1200068.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1016 Peak 1 | 0.0130<br>0.0104 | 0.0132 | 0.0125 | 0.0117 | 0.0109 | Lin1 | 0.0002 | 0.0106 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1016 Peak 2 | 0.0226<br>0.0190 | 0.0238 | 0.0222 | 0.0210 | 0.0197 | Lin1 | 0.0003 | 0.0193 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1016 Peak 3 | 0.0471<br>0.0469 | 0.0503 | 0.0465 | 0.0479 | 0.0469 | Lin1 | 0.0001 | 0.0469 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 4 | 0.0212<br>0.0189 | 0.0222 | 0.0202 | 0.0199 | 0.0193 | Lin1 | 0.0002 | 0.0190 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 5 | 0.0220<br>0.0187 | 0.0228 | 0.0210 | 0.0199 | 0.0191 | Lin1 | 0.0003 | 0.0188 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1260 Peak 1 | 0.0330<br>0.0301 | 0.0338 | 0.0324 | 0.0309 | 0.0302 | Lin1 | 0.0003 | 0.0300 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 2 | 0.0622<br>0.0585 | 0.0618 | 0.0622 | 0.0606 | 0.0586 | Lin1 | 0.0003 | 0.0586 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 3 | 0.0650<br>0.0628 | 0.0634 | 0.0652 | 0.0643 | 0.0626 | Lin1 | 0.0002 | 0.0629 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 4 | 0.0661<br>0.0714 | 0.0711 | 0.0771 | 0.0773 | 0.0731 | Lin1 | 0 | 0.0729 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1260 Peak 5 | 0.0374<br>0.0384 | 0.0399 | 0.0382 | 0.0386 | 0.0385 | Lin1 | 0 | 0.0384 | | | | | | | 1.0000 | | 0.9900 |
| Tetrachloro-m-xylene | 0.7048<br>0.6608 | 0.7003 | 0.7144 | 0.6986 | 0.6822 | Lin1 | 0.0002 | 0.6699 | | | | | | | 0.9990 | | 0.9900 |
| DCB Decachlorobiphenyl | 0.7699<br>0.6896 | 0.7629 | 0.7446 | 0.7316 | 0.6953 | Lin1 | 0.0003 | 0.6927 | | | | | | | 1.0000 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-57930-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/17/2015 03:29 Calibration End Date: 11/17/2015 04:59 Calibration ID: 31555

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/28 | P1200063.D |
| Level 2 | STD01 240-207007/29 | P1200064.D |
| Level 3 | STD02 240-207007/30 | P1200065.D |
| Level 4 | STD05 240-207007/31 | P1200066.D |
| Level 5 | STD1 240-207007/32 | P1200067.D |
| Level 6 | STD2 240-207007/33 | P1200068.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1016 Peak 1 | BNB | Lin1 | 2707871<br>80765612 | 4820722 | 9085846 | 22006625 | 42447741 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 2 | BNB | Lin1 | 4700873<br>147609677 | 8734684 | 16183493 | 39469767 | 76715703 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 3 | BNB | Lin1 | 9792259<br>364184674 | 18438138 | 33886380 | 90045825 | 182657834 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 4 | BNB | Lin1 | 4396897<br>147122332 | 8128551 | 14710049 | 37512861 | 75171974 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 5 | BNB | Lin1 | 4563462<br>145010499 | 8352402 | 15341259 | 37426880 | 74524891 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 1 | BNB | Lin1 | 6849336<br>233433821 | 12383500 | 23597918 | 58080604 | 117631104 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 2 | BNB | Lin1 | 12922388<br>454390240 | 22631342 | 45372536 | 113933482 | 228581266 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 3 | BNB | Lin1 | 13513426<br>488067042 | 23245320 | 47522931 | 120987145 | 244049421 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 4 | BNB | Lin1 | 13737456<br>554191846 | 26050869 | 56208655 | 145378951 | 285103725 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 5 | BNB | Lin1 | 7764692<br>298012472 | 14617316 | 27852633 | 72613053 | 149980149 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| Tetrachloro-m-xylene | BNB | Lin1 | 7322625<br>256611202 | 12827811 | 26051804 | 65705303 | 132973208 | 0.00250<br>0.100 | 0.00500 | 0.0100 | 0.0250 | 0.0500 |
| DCB Decachlorobiphenyl | BNB | Lin1 | 7998062<br>267771950 | 13974803 | 27151934 | 68816182 | 135510483 | 0.00250<br>0.100 | 0.00500 | 0.0100 | 0.0250 | 0.0500 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-207007/39  Calibration Date: 11/17/2015 06:48

Instrument ID: A2HP12  Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/17/2015 04:59

Lab File ID: P1200074.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0109 | | 1.01 | 1.00 | 1.2 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0209 | | 1.07 | 1.00 | 6.9 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0504 | | 1.07 | 1.00 | 7.2 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0205 | | 1.07 | 1.00 | 6.9 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0196 | | 1.03 | 1.00 | 3.0 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0302 | | 0.995 | 1.00 | -0.5 | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0570 | | 0.967 | 1.00 | -3.3 | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0535 | | 0.848 | 1.00 | -15.2 | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0824 | | 1.13 | 1.00 | 13.1 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0405 | | 1.05 | 1.00 | 5.3 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6836 | | 0.0507 | 0.0500 | 1.5 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.6921 | | 0.0495 | 0.0500 | -1.0 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: ICV 240-207007/39 Calibration Date: 11/17/2015 06:48

Instrument ID: A2HP12 Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 04:59

Lab File ID: P1200074.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1016 Peak 1 | 4.61 | 4.58 | 4.64 |
| PCB-1016 Peak 2 | 5.11 | 5.08 | 5.14 |
| PCB-1016 Peak 3 | 5.74 | 5.71 | 5.77 |
| PCB-1016 Peak 4 | 5.91 | 5.88 | 5.94 |
| PCB-1016 Peak 5 | 6.45 | 6.42 | 6.48 |
| PCB-1260 Peak 1 | 7.95 | 7.92 | 7.98 |
| PCB-1260 Peak 2 | 8.29 | 8.26 | 8.32 |
| PCB-1260 Peak 3 | 8.61 | 8.58 | 8.64 |
| PCB-1260 Peak 4 | 9.40 | 9.37 | 9.43 |
| PCB-1260 Peak 5 | 9.72 | 9.69 | 9.75 |
| Tetrachloro-m-xylene | 4.05 | 4.02 | 4.08 |
| DCB Decachlorobiphenyl | 10.93 | 10.90 | 10.96 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/3  Calibration Date: 11/18/2015 18:06

Instrument ID: A2HP12  Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/17/2015 04:59

Lab File ID: P1200039.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0113 | | 0.513 | 0.500 | 2.5 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0196 | | 0.490 | 0.500 | -1.9 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0438 | | 0.465 | 0.500 | -7.1 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0181 | | 0.466 | 0.500 | -6.7 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0184 | | 0.476 | 0.500 | -4.9 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0286 | | 0.467 | 0.500 | -6.7 | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0526 | | 0.443 | 0.500 | -11.4 | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0569 | | 0.450 | 0.500 | -10.0 | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0644 | | 0.442 | 0.500 | -11.6 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0336 | | 0.437 | 0.500 | -12.7 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6747 | | 0.0249 | 0.0250 | -0.4 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.5804 | | 0.0205 | 0.0250 | -18.0 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/3 Calibration Date: 11/18/2015 18:06

Instrument ID: A2HP12 Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 04:59

Lab File ID: P1200039.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1016 Peak 1 | 4.61 | 4.58 | 4.64 |
| PCB-1016 Peak 2 | 5.11 | 5.08 | 5.14 |
| PCB-1016 Peak 3 | 5.74 | 5.71 | 5.77 |
| PCB-1016 Peak 4 | 5.91 | 5.88 | 5.94 |
| PCB-1016 Peak 5 | 6.45 | 6.42 | 6.48 |
| PCB-1260 Peak 1 | 7.95 | 7.92 | 7.98 |
| PCB-1260 Peak 2 | 8.29 | 8.26 | 8.32 |
| PCB-1260 Peak 3 | 8.61 | 8.58 | 8.64 |
| PCB-1260 Peak 4 | 9.40 | 9.37 | 9.43 |
| PCB-1260 Peak 5 | 9.72 | 9.69 | 9.75 |
| Tetrachloro-m-xylene | 4.05 | 4.02 | 4.08 |
| DCB Decachlorobiphenyl | 10.93 | 10.90 | 10.96 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/4  Calibration Date: 11/18/2015 18:24

Instrument ID: A2HP12  Calib Start Date: 11/16/2015 20:14

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/16/2015 21:45

Lab File ID: P1200040.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1232 Peak 1 | Lin1 | | 0.0135 | | 0.536 | 0.500 | 7.3 | 20.0 |
| PCB-1232 Peak 2 | Lin1 | | 0.0095 | | 0.519 | 0.500 | 3.7 | 20.0 |
| PCB-1232 Peak 3 | Lin1 | | 0.0210 | | 0.504 | 0.500 | 0.7 | 20.0 |
| PCB-1232 Peak 4 | Lin1 | | 0.0090 | | 0.528 | 0.500 | 5.7 | 20.0 |
| PCB-1232 Peak 5 | Lin1 | | 0.0082 | | 0.517 | 0.500 | 3.4 | 20.0 |
| PCB-1262 Peak 1 | Lin1 | | 0.0240 | | 0.506 | 0.500 | 1.2 | 20.0 |
| PCB-1262 Peak 2 | Lin1 | | 0.0409 | | 0.456 | 0.500 | -8.7 | 20.0 |
| PCB-1262 Peak 3 | Lin1 | | 0.0374 | | 0.455 | 0.500 | -9.1 | 20.0 |
| PCB-1262 Peak 4 | Lin1 | | 0.0933 | | 0.482 | 0.500 | -3.6 | 20.0 |
| PCB-1262 Peak 5 | Lin1 | | 0.0356 | | 0.466 | 0.500 | -6.8 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/4 Calibration Date: 11/18/2015 18:24

Instrument ID: A2HP12 Calib Start Date: 11/16/2015 20:14

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/16/2015 21:45

Lab File ID: P1200040.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1232 Peak 1 | 4.61 | 4.58 | 4.64 |
| PCB-1232 Peak 2 | 5.11 | 5.08 | 5.14 |
| PCB-1232 Peak 3 | 5.74 | 5.71 | 5.77 |
| PCB-1232 Peak 4 | 5.91 | 5.88 | 5.94 |
| PCB-1232 Peak 5 | 6.45 | 6.42 | 6.48 |
| PCB-1262 Peak 1 | 7.96 | 7.93 | 7.99 |
| PCB-1262 Peak 2 | 8.75 | 8.72 | 8.78 |
| PCB-1262 Peak 3 | 9.03 | 9.00 | 9.06 |
| PCB-1262 Peak 4 | 9.40 | 9.37 | 9.43 |
| PCB-1262 Peak 5 | 9.77 | 9.74 | 9.80 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/5 Calibration Date: 11/18/2015 18:43

Instrument ID: A2HP12 Calib Start Date: 11/16/2015 22:03

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/16/2015 23:33

Lab File ID: P1200041.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1242 Peak 1 | Lin1 | | 0.0089 | | 0.516 | 0.500 | 3.3 | 20.0 |
| PCB-1242 Peak 2 | Lin1 | | 0.0161 | | 0.501 | 0.500 | 0.2 | 20.0 |
| PCB-1242 Peak 3 | Lin1 | | 0.0361 | | 0.477 | 0.500 | -4.5 | 20.0 |
| PCB-1242 Peak 4 | Lin1 | | 0.0154 | | 0.493 | 0.500 | -1.3 | 20.0 |
| PCB-1242 Peak 5 | Lin1 | | 0.0156 | | 0.490 | 0.500 | -2.1 | 20.0 |
| PCB-1268 Peak 1 | Lin1 | | 0.0979 | | 0.438 | 0.500 | -12.5 | 20.0 |
| PCB-1268 Peak 2 | Lin1 | | 0.0977 | | 0.449 | 0.500 | -10.3 | 20.0 |
| PCB-1268 Peak 3 | Lin1 | | 0.0830 | | 0.448 | 0.500 | -10.4 | 20.0 |
| PCB-1268 Peak 4 | Lin1 | | 0.0338 | | 0.441 | 0.500 | -11.8 | 20.0 |
| PCB-1268 Peak 5 | Lin1 | | 0.2750 | | 0.438 | 0.500 | -12.4 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/5 Calibration Date: 11/18/2015 18:43

Instrument ID: A2HP12 Calib Start Date: 11/16/2015 22:03

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/16/2015 23:33

Lab File ID: P1200041.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1242 Peak 1 | 4.61 | 4.58 | 4.64 |
| PCB-1242 Peak 2 | 5.11 | 5.08 | 5.14 |
| PCB-1242 Peak 3 | 5.74 | 5.71 | 5.77 |
| PCB-1242 Peak 4 | 5.91 | 5.88 | 5.94 |
| PCB-1242 Peak 5 | 6.45 | 6.42 | 6.48 |
| PCB-1268 Peak 1 | 9.77 | 9.74 | 9.80 |
| PCB-1268 Peak 2 | 9.81 | 9.78 | 9.84 |
| PCB-1268 Peak 3 | 10.03 | 10.00 | 10.06 |
| PCB-1268 Peak 4 | 10.44 | 10.41 | 10.47 |
| PCB-1268 Peak 5 | 10.74 | 10.71 | 10.77 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/6  Calibration Date: 11/18/2015 19:01

Instrument ID: A2HP12  Calib Start Date: 11/16/2015 23:51

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/17/2015 01:22

Lab File ID: P1200042.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1248 Peak 1 | Lin1 | | 0.0088 | | 0.508 | 0.500 | 1.7 | 20.0 |
| PCB-1248 Peak 2 | Lin1 | | 0.0233 | | 0.478 | 0.500 | -4.3 | 20.0 |
| PCB-1248 Peak 3 | Lin1 | | 0.0215 | | 0.478 | 0.500 | -4.4 | 20.0 |
| PCB-1248 Peak 4 | Lin1 | | 0.0183 | | 0.475 | 0.500 | -5.0 | 20.0 |
| PCB-1248 Peak 5 | Lin1 | | 0.0121 | | 0.486 | 0.500 | -2.8 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/6 Calibration Date: 11/18/2015 19:01

Instrument ID: A2HP12 Calib Start Date: 11/16/2015 23:51

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 01:22

Lab File ID: P1200042.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1248 Peak 1 | 5.11 | 5.08 | 5.14 |
| PCB-1248 Peak 2 | 5.74 | 5.71 | 5.77 |
| PCB-1248 Peak 3 | 7.01 | 6.98 | 7.04 |
| PCB-1248 Peak 4 | 7.26 | 7.23 | 7.29 |
| PCB-1248 Peak 5 | 7.73 | 7.70 | 7.76 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/7  Calibration Date: 11/18/2015 19:19

Instrument ID: A2HP12  Calib Start Date: 11/17/2015 01:40

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/17/2015 03:10

Lab File ID: P1200043.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1221 Peak 1 | Lin1 | | 0.0065 | | 0.535 | 0.500 | 7.0 | 20.0 |
| PCB-1221 Peak 2 | Lin1 | | 0.0043 | | 0.543 | 0.500 | 8.6 | 20.0 |
| PCB-1221 Peak 3 | Lin1 | | 0.0165 | | 0.536 | 0.500 | 7.1 | 20.0 |
| PCB-1254 Peak 1 | Lin1 | | 0.0213 | | 0.517 | 0.500 | 3.5 | 20.0 |
| PCB-1254 Peak 2 | Lin1 | | 0.0299 | | 0.508 | 0.500 | 1.6 | 20.0 |
| PCB-1254 Peak 3 | Lin1 | | 0.0445 | | 0.537 | 0.500 | 7.5 | 20.0 |
| PCB-1254 Peak 4 | Lin1 | | 0.0372 | | 0.516 | 0.500 | 3.2 | 20.0 |
| PCB-1254 Peak 5 | Lin1 | | 0.0463 | | 0.502 | 0.500 | 0.3 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCV 240-207364/7 Calibration Date: 11/18/2015 19:19

Instrument ID: A2HP12 Calib Start Date: 11/17/2015 01:40

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 03:10

Lab File ID: P1200043.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1221 Peak 1 | 4.34 | 4.31 | 4.37 |
| PCB-1221 Peak 2 | 4.55 | 4.52 | 4.58 |
| PCB-1221 Peak 3 | 4.61 | 4.58 | 4.64 |
| PCB-1254 Peak 1 | 6.93 | 6.90 | 6.96 |
| PCB-1254 Peak 2 | 7.26 | 7.23 | 7.29 |
| PCB-1254 Peak 3 | 7.73 | 7.70 | 7.76 |
| PCB-1254 Peak 4 | 8.07 | 8.04 | 8.10 |
| PCB-1254 Peak 5 | 8.62 | 8.59 | 8.65 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207364/28  Calibration Date: 11/19/2015 01:40

Instrument ID: A2HP12  Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/17/2015 04:59

Lab File ID: P1200064.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0112 | | 0.508 | 0.500 | 1.6 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0198 | | 0.495 | 0.500 | -0.9 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0450 | | 0.478 | 0.500 | -4.5 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0186 | | 0.477 | 0.500 | -4.6 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0187 | | 0.482 | 0.500 | -3.7 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0291 | | 0.476 | 0.500 | -4.9 | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0555 | | 0.468 | 0.500 | -6.4 | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0589 | | 0.466 | 0.500 | -6.9 | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0652 | | 0.448 | 0.500 | -10.4 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0349 | | 0.453 | 0.500 | -9.4 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6633 | | 0.0245 | 0.0250 | -2.1 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.5855 | | 0.0207 | 0.0250 | -17.3 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Lab Sample ID: CCVIS 240-207364/28 Calibration Date: 11/19/2015 01:40

Instrument ID: A2HP12 Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 04:59

Lab File ID: P1200064.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1016 Peak 1 | 4.61 | 4.58 | 4.64 |
| PCB-1016 Peak 2 | 5.11 | 5.08 | 5.14 |
| PCB-1016 Peak 3 | 5.74 | 5.71 | 5.77 |
| PCB-1016 Peak 4 | 5.91 | 5.88 | 5.94 |
| PCB-1016 Peak 5 | 6.45 | 6.42 | 6.48 |
| PCB-1260 Peak 1 | 7.96 | 7.93 | 7.99 |
| PCB-1260 Peak 2 | 8.29 | 8.26 | 8.32 |
| PCB-1260 Peak 3 | 8.61 | 8.58 | 8.64 |
| PCB-1260 Peak 4 | 9.40 | 9.37 | 9.43 |
| PCB-1260 Peak 5 | 9.72 | 9.69 | 9.75 |
| Tetrachloro-m-xylene | 4.05 | 4.02 | 4.08 |
| DCB Decachlorobiphenyl | 10.93 | 10.90 | 10.96 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206917/21-A

Matrix: Solid | Lab File ID: P1200062.D

Analysis Method: 8082A | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 01:03

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 12 | U | 33 | 12 |
| 11104-28-2 | Aroclor-1221 | 16 | U | 33 | 16 |
| 11141-16-5 | Aroclor-1232 | 20 | U | 33 | 20 |
| 53469-21-9 | Aroclor-1242 | 11 | U | 33 | 11 |
| 12672-29-6 | Aroclor-1248 | 8.0 | U | 33 | 8.0 |
| 11097-69-1 | Aroclor-1254 | 14 | U | 33 | 14 |
| 11096-82-5 | Aroclor-1260 | 9.0 | U | 33 | 9.0 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 74 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 79 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-206917/22-A

Matrix: Solid | Lab File ID: P1200063.D

Analysis Method: 8082A | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 30(g) | Date Analyzed: 11/19/2015 01:21

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 251 | | 33 | 12 |
| 11096-82-5 | Aroclor-1260 | 253 | | 33 | 9.0 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 75 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 77 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57880-E-1-B MS

Matrix: Solid | Lab File ID: P1200066.D

Analysis Method: 8082A | Date Collected: 11/12/2015 10:30

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 29.95(g) | Date Analyzed: 11/19/2015 02:16

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 24.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 518 | | 44 | 16 |
| 11104-28-2 | Aroclor-1221 | 21 | U | 44 | 21 |
| 11141-16-5 | Aroclor-1232 | 27 | U | 44 | 27 |
| 53469-21-9 | Aroclor-1242 | 15 | U | 44 | 15 |
| 12672-29-6 | Aroclor-1248 | 11 | U | 44 | 11 |
| 11097-69-1 | Aroclor-1254 | 19 | U | 44 | 19 |
| 11096-82-5 | Aroclor-1260 | 266 | | 44 | 12 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 75 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 94 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57930-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-57880-E-1-C MSD

Matrix: Solid | Lab File ID: P1200067.D

Analysis Method: 8082A | Date Collected: 11/12/2015 10:30

Extraction Method: 3540C | Date Extracted: 11/16/2015 09:32

Sample wt/vol: 29.75(g) | Date Analyzed: 11/19/2015 02:34

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 24.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 207364 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 638 | | 44 | 16 |
| 11104-28-2 | Aroclor-1221 | 21 | U | 44 | 21 |
| 11141-16-5 | Aroclor-1232 | 27 | U | 44 | 27 |
| 53469-21-9 | Aroclor-1242 | 15 | U | 44 | 15 |
| 12672-29-6 | Aroclor-1248 | 11 | U | 44 | 11 |
| 11097-69-1 | Aroclor-1254 | 19 | U | 44 | 19 |
| 11096-82-5 | Aroclor-1260 | 331 | | 44 | 12 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 90 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 87 | | 10-155 |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP12 Start Date: 11/16/2015 20:14

Analysis Batch Number: 207007 End Date: 11/17/2015 07:24

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD005 240-207007/4 IC | | 11/16/2015 20:14 | 1 | P1200039.D | CLP-1 0.53(mm) |
| STD005 240-207007/4 IC | | 11/16/2015 20:14 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/5 IC | | 11/16/2015 20:32 | 1 | P1200040.D | CLP-1 0.53(mm) |
| STD01 240-207007/5 IC | | 11/16/2015 20:32 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/6 IC | | 11/16/2015 20:50 | 1 | P1200041.D | CLP-1 0.53(mm) |
| STD02 240-207007/6 IC | | 11/16/2015 20:50 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/7 IC | | 11/16/2015 21:09 | 1 | P1200042.D | CLP-1 0.53(mm) |
| STD05 240-207007/7 IC | | 11/16/2015 21:09 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/8 IC | | 11/16/2015 21:27 | 1 | P1200043.D | CLP-1 0.53(mm) |
| STD1 240-207007/8 IC | | 11/16/2015 21:27 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/9 IC | | 11/16/2015 21:45 | 1 | P1200044.D | CLP-1 0.53(mm) |
| STD2 240-207007/9 IC | | 11/16/2015 21:45 | 1 | | CLP-2 0.53(mm) |
| STD005 240-207007/10 IC | | 11/16/2015 22:03 | 1 | P1200045.D | CLP-1 0.53(mm) |
| STD005 240-207007/10 IC | | 11/16/2015 22:03 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/11 IC | | 11/16/2015 22:21 | 1 | P1200046.D | CLP-1 0.53(mm) |
| STD01 240-207007/11 IC | | 11/16/2015 22:21 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/12 IC | | 11/16/2015 22:39 | 1 | P1200047.D | CLP-1 0.53(mm) |
| STD02 240-207007/12 IC | | 11/16/2015 22:39 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/13 IC | | 11/16/2015 22:57 | 1 | P1200048.D | CLP-1 0.53(mm) |
| STD05 240-207007/13 IC | | 11/16/2015 22:57 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/14 IC | | 11/16/2015 23:15 | 1 | P1200049.D | CLP-1 0.53(mm) |
| STD1 240-207007/14 IC | | 11/16/2015 23:15 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/15 IC | | 11/16/2015 23:33 | 1 | P1200050.D | CLP-1 0.53(mm) |
| STD2 240-207007/15 IC | | 11/16/2015 23:33 | 1 | | CLP-2 0.53(mm) |
| STD005 240-207007/16 IC | | 11/16/2015 23:51 | 1 | P1200051.D | CLP-1 0.53(mm) |
| STD005 240-207007/16 IC | | 11/16/2015 23:51 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/17 IC | | 11/17/2015 00:09 | 1 | P1200052.D | CLP-1 0.53(mm) |
| STD01 240-207007/17 IC | | 11/17/2015 00:09 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/18 IC | | 11/17/2015 00:28 | 1 | P1200053.D | CLP-1 0.53(mm) |
| STD02 240-207007/18 IC | | 11/17/2015 00:28 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/19 IC | | 11/17/2015 00:46 | 1 | P1200054.D | CLP-1 0.53(mm) |
| STD05 240-207007/19 IC | | 11/17/2015 00:46 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/20 IC | | 11/17/2015 01:04 | 1 | P1200055.D | CLP-1 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP12 Start Date: 11/16/2015 20:14

Analysis Batch Number: 207007 End Date: 11/17/2015 07:24

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-207007/20 IC | | 11/17/2015 01:04 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/21 IC | | 11/17/2015 01:22 | 1 | P1200056.D | CLP-1 0.53(mm) |
| STD2 240-207007/21 IC | | 11/17/2015 01:22 | 1 | | CLP-2 0.53(mm) |
| STD005 240-207007/22 IC | | 11/17/2015 01:40 | 1 | P1200057.D | CLP-1 0.53(mm) |
| STD005 240-207007/22 IC | | 11/17/2015 01:40 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/23 IC | | 11/17/2015 01:58 | 1 | P1200058.D | CLP-1 0.53(mm) |
| STD01 240-207007/23 IC | | 11/17/2015 01:58 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/24 IC | | 11/17/2015 02:16 | 1 | P1200059.D | CLP-1 0.53(mm) |
| STD02 240-207007/24 IC | | 11/17/2015 02:16 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/25 IC | | 11/17/2015 02:34 | 1 | P1200060.D | CLP-1 0.53(mm) |
| STD05 240-207007/25 IC | | 11/17/2015 02:34 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/26 IC | | 11/17/2015 02:52 | 1 | P1200061.D | CLP-1 0.53(mm) |
| STD1 240-207007/26 IC | | 11/17/2015 02:52 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/27 IC | | 11/17/2015 03:10 | 1 | P1200062.D | CLP-1 0.53(mm) |
| STD2 240-207007/27 IC | | 11/17/2015 03:10 | 1 | | CLP-2 0.53(mm) |
| STD005 240-207007/28 IC | | 11/17/2015 03:29 | 1 | P1200063.D | CLP-1 0.53(mm) |
| STD005 240-207007/28 IC | | 11/17/2015 03:29 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/29 IC | | 11/17/2015 03:47 | 1 | P1200064.D | CLP-1 0.53(mm) |
| STD01 240-207007/29 IC | | 11/17/2015 03:47 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/30 IC | | 11/17/2015 04:05 | 1 | P1200065.D | CLP-1 0.53(mm) |
| STD02 240-207007/30 IC | | 11/17/2015 04:05 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/31 ICIS | | 11/17/2015 04:23 | 1 | P1200066.D | CLP-1 0.53(mm) |
| STD05 240-207007/31 ICIS | | 11/17/2015 04:23 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/32 IC | | 11/17/2015 04:41 | 1 | P1200067.D | CLP-1 0.53(mm) |
| STD1 240-207007/32 IC | | 11/17/2015 04:41 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/33 IC | | 11/17/2015 04:59 | 1 | P1200068.D | CLP-1 0.53(mm) |
| STD2 240-207007/33 IC | | 11/17/2015 04:59 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/34 | | 11/17/2015 05:17 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/34 | | 11/17/2015 05:17 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/35 | | 11/17/2015 05:35 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/35 | | 11/17/2015 05:35 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/36 | | 11/17/2015 05:53 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/36 | | 11/17/2015 05:53 | 1 | | CLP-2 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP12  Start Date: 11/16/2015 20:14

Analysis Batch Number: 207007  End Date: 11/17/2015 07:24

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| ICV 240-207007/37 | | 11/17/2015 06:12 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/37 | | 11/17/2015 06:12 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/38 | | 11/17/2015 06:30 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/38 | | 11/17/2015 06:30 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/39 | | 11/17/2015 06:48 | 1 | P1200074.D | CLP-1 0.53(mm) |
| ICV 240-207007/39 | | 11/17/2015 06:48 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/40 | | 11/17/2015 07:06 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/40 | | 11/17/2015 07:06 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/41 | | 11/17/2015 07:24 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/41 | | 11/17/2015 07:24 | 1 | | CLP-2 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP12 Start Date: 11/18/2015 18:06

Analysis Batch Number: 207364 End Date: 11/19/2015 02:34

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-207364/3 CCVIS | | 11/18/2015 18:06 | 1 | P1200039.D | CLP-1 0.53(mm) |
| CCV 240-207364/3 CCVIS | | 11/18/2015 18:06 | 1 | | CLP-2 0.53(mm) |
| CCV 240-207364/4 | | 11/18/2015 18:24 | 1 | P1200040.D | CLP-1 0.53(mm) |
| CCV 240-207364/4 | | 11/18/2015 18:24 | 1 | | CLP-2 0.53(mm) |
| CCV 240-207364/5 | | 11/18/2015 18:43 | 1 | P1200041.D | CLP-1 0.53(mm) |
| CCV 240-207364/5 | | 11/18/2015 18:43 | 1 | | CLP-2 0.53(mm) |
| CCV 240-207364/6 | | 11/18/2015 19:01 | 1 | P1200042.D | CLP-1 0.53(mm) |
| CCV 240-207364/6 | | 11/18/2015 19:01 | 1 | | CLP-2 0.53(mm) |
| CCV 240-207364/7 | | 11/18/2015 19:19 | 1 | P1200043.D | CLP-1 0.53(mm) |
| CCV 240-207364/7 | | 11/18/2015 19:19 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 11/18/2015 19:55 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 19:55 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 11/18/2015 20:13 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 20:13 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 11/18/2015 21:44 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 21:44 | 1 | | CLP-2 0.53(mm) |
| 240-57930-6 | F-VB-UST-P-2 | 11/18/2015 22:02 | 1 | P1200052.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 22:02 | 1 | | CLP-2 0.53(mm) |
| 240-57930-7 | F-VB-UST-P-3 | 11/18/2015 22:20 | 1 | P1200053.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 22:20 | 1 | | CLP-2 0.53(mm) |
| 240-57930-8 | F-VB-UST-P-4 | 11/18/2015 22:38 | 1 | P1200054.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 22:38 | 1 | | CLP-2 0.53(mm) |
| 240-57930-9 | F-VB-UST-P-5 | 11/18/2015 22:56 | 1 | P1200055.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 22:56 | 1 | | CLP-2 0.53(mm) |
| 240-57930-10 | F-VB-UST-P-6 | 11/18/2015 23:15 | 1 | P1200056.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 23:15 | 1 | | CLP-2 0.53(mm) |
| 240-57930-11 | F-VS-UST-P-DUP | 11/18/2015 23:33 | 1 | P1200057.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 23:33 | 1 | | CLP-2 0.53(mm) |
| 240-57930-1 | F-VS-UST-P-1 | 11/18/2015 23:51 | 1 | P1200058.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/18/2015 23:51 | 1 | | CLP-2 0.53(mm) |
| 240-57930-2 | F-VS-UST-P-2 | 11/19/2015 00:09 | 1 | P1200059.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/19/2015 00:09 | 1 | | CLP-2 0.53(mm) |
| 240-57930-3 | F-VS-UST-P-3 | 11/19/2015 00:27 | 1 | P1200060.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/19/2015 00:27 | 1 | | CLP-2 0.53(mm) |
| 240-57930-4 | F-VS-UST-P-4 | 11/19/2015 00:45 | 1 | P1200061.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/19/2015 00:45 | 1 | | CLP-2 0.53(mm) |
| MB 240-206917/21-A | | 11/19/2015 01:03 | 1 | P1200062.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/19/2015 01:03 | 1 | | CLP-2 0.53(mm) |
| LCS 240-206917/22-A | | 11/19/2015 01:21 | 1 | P1200063.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/19/2015 01:21 | 1 | | CLP-2 0.53(mm) |
| CCVIS 240-207364/28 | | 11/19/2015 01:40 | 1 | P1200064.D | CLP-1 0.53(mm) |
| CCVIS 240-207364/28 | | 11/19/2015 01:40 | 1 | | CLP-2 0.53(mm) |
| 240-57930-5 | F-VB-UST-P-1 | 11/19/2015 01:58 | 1 | P1200065.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/19/2015 01:58 | 1 | | CLP-2 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Instrument ID: A2HP12  Start Date: 11/18/2015 18:06

Analysis Batch Number: 207364  End Date: 11/19/2015 02:34

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| 240-57880-E-1-B MS | | 11/19/2015 02:16 | 1 | P1200066.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/19/2015 02:16 | 1 | | CLP-2 0.53(mm) |
| 240-57880-E-1-C MSD | | 11/19/2015 02:34 | 1 | P1200067.D | CLP-1 0.53(mm) |
| ZZZZZ | | 11/19/2015 02:34 | 1 | | CLP-2 0.53(mm) |

# PCBS BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 206917  Batch Start Date: 11/16/15 09:32  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | ex10PPMSPK 00018 | ex2/.2SURRW 00062 | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57930-D-6 | F-VB-UST-P-2 | 3540C, 8082A | T | 29.88 g | 10 mL | | 1 mL | | |
| 240-57930-D-7 | F-VB-UST-P-3 | 3540C, 8082A | T | 30.01 g | 10 mL | | 1 mL | | |
| 240-57930-D-8 | F-VB-UST-P-4 | 3540C, 8082A | T | 29.74 g | 10 mL | | 1 mL | | |
| 240-57930-D-9 | F-VB-UST-P-5 | 3540C, 8082A | T | 29.69 g | 10 mL | | 1 mL | | |
| 240-57930-D-10 | F-VB-UST-P-6 | 3540C, 8082A | T | 29.68 g | 10 mL | | 1 mL | | |
| 240-57930-D-11 | F-VS-UST-P-DUP | 3540C, 8082A | T | 29.83 g | 10 mL | | 1 mL | | |
| 240-57930-D-1 | F-VS-UST-P-1 | 3540C, 8082A | T | 29.62 g | 10 mL | | 1 mL | | |
| 240-57930-D-2 | F-VS-UST-P-2 | 3540C, 8082A | T | 30.44 g | 10 mL | | 1 mL | | |
| 240-57930-D-3 | F-VS-UST-P-3 | 3540C, 8082A | T | 29.65 g | 10 mL | | 1 mL | | |
| 240-57930-D-4 | F-VS-UST-P-4 | 3540C, 8082A | T | 30.41 g | 10 mL | | 1 mL | | |
| 240-57930-D-5 | F-VB-UST-P-1 | 3540C, 8082A | T | 30.45 g | 10 mL | | 1 mL | | |
| 240-57880-E-1 MS | | 3540C, 8082A | T | 29.95 g | 10 mL | 1 mL | 1 mL | | |
| 240-57880-E-1 MSD | | 3540C, 8082A | T | 29.75 g | 10 mL | 1 mL | 1 mL | | |
| MB 240-206917/21 | | 3540C, 8082A | | 30 g | 10 mL | | 1 mL | | |
| LCS 240-206917/22 | | 3540C, 8082A | | 30 g | 10 mL | 1 mL | 1 mL | | |

| Batch Notes | |
|---|---|
| Acid used for Clean Up Reagent | 2226516 |
| Balance ID | 1339907 |
| Na2SO4 Lot Number | 2204509 |
| Prep Solvent Lot # | 2286385 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GENERAL CHEMISTRY

COVER PAGE
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton  Job Number: 240-57930-1

SDG No.:

Project: LM MRC Block F UST Pipe Inv.

| Client Sample ID | Lab Sample ID |
|---|---|
| F-VS-UST-P-1 | 240-57930-1 |
| F-VS-UST-P-2 | 240-57930-2 |
| F-VS-UST-P-3 | 240-57930-3 |
| F-VS-UST-P-4 | 240-57930-4 |
| F-VB-UST-P-1 | 240-57930-5 |
| F-VB-UST-P-2 | 240-57930-6 |
| F-VB-UST-P-3 | 240-57930-7 |
| F-VB-UST-P-4 | 240-57930-8 |
| F-VB-UST-P-5 | 240-57930-9 |
| F-VB-UST-P-6 | 240-57930-10 |
| F-VS-UST-P-DUP | 240-57930-11 |

Comments:

9-IN
DETECTION LIMITS
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job Number: 240-57930-1

SDG Number:

Matrix: Solid Instrument ID: NOEQUIP

Method: Moisture RL Date: 01/28/2010 09:24

| Analyte | Wavelength/ Mass | RL (%) | |
|---|---|---|---|
| Percent Moisture | | 0.1 | |
| Percent Solids | | 0.1 | |

9-IN
CALIBRATION BLANK DETECTION LIMITS
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton  Job Number: 240-57930-1

SDG Number:

Matrix: Solid  Instrument ID: NOEQUIP

Method: Moisture  XRL Date: 01/28/2010 09:24

| Analyte | Wavelength/ Mass | XRL (mg/L) | |
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| Percent Moisture | | 10 | |
| Percent Solids | | 10 | |

13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: NOEQUIP Method: Moisture

Start Date: 11/16/2015 15:41 End Date: 11/16/2015 16:19

| Lab Sample ID | D / F | Type | Time | Analytes | | | | | | | | | | | | | | | | | | | | |
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| 240-57930-11 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-10 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-9 DU | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-9 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-8 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-7 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-6 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-5 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-4 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-3 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-2 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
| 240-57930-1 | 1 | T | 15:41 | X | X | | | | | | | | | | | | | | | | | | | |
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13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: NOEQUIP Method: Moisture

Start Date: 11/16/2015 15:41 End Date: 11/16/2015 16:19

| Lab Sample ID | D / F | Type | Time | Analytes | | | | | | | | | | | | | | | | | | | |
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13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: NOEQUIP Method: Moisture

Start Date: 11/16/2015 15:41 End Date: 11/16/2015 16:19

| Lab Sample ID | D / F | Type | Time | Analytes | | | | | | | | | | | | | | | | | | | |
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13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: NOEQUIP Method: Moisture

Start Date: 11/16/2015 15:41 End Date: 11/16/2015 16:19

| Lab Sample ID | D / F | Type | Time | Analytes | | | | | | | | | | | | | | | | | | | |
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13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job No.: 240-57930-1

SDG No.:

Instrument ID: NOEQUIP Method: Moisture

Start Date: 11/16/2015 15:41 End Date: 11/16/2015 16:19

| Lab Sample ID | D / F | Type | Time | Analytes | | | | | | | | | | | | | | | | | | | | |
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Prep Types

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# GENERAL CHEMISTRY BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57930-1

SDG No.:

Batch Number: 207016  Batch Start Date: 11/16/15 15:41  Batch Analyst: Nguyen, Diem

Batch Method: Moisture  Batch End Date: 11/17/15 11:10

| Lab Sample ID | Client Sample ID | Method Chain | Basis | DishWeight | SampleMassWet | SampleMassDry | | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57930-A-11 | F-VS-UST-P-DUP | Moisture | T | 4.3999 g | 25.3910 g | 22.1197 g | | | |
| 240-57930-A-10 | F-VB-UST-P-6 | Moisture | T | 4.3999 g | 20.4530 g | 18.0088 g | | | |
| 240-57930-A-9 DU | F-VB-UST-P-5 | Moisture | T | 4.3999 g | 12.2743 g | 10.9709 g | | | |
| 240-57930-A-9 | F-VB-UST-P-5 | Moisture | T | 4.3999 g | 13.5764 g | 12.0435 g | | | |
| 240-57930-A-8 | F-VB-UST-P-4 | Moisture | T | 4.3999 g | 19.2510 g | 16.8544 g | | | |
| 240-57930-A-7 | F-VB-UST-P-3 | Moisture | T | 4.3999 g | 19.6277 g | 16.7311 g | | | |
| 240-57930-A-6 | F-VB-UST-P-2 | Moisture | T | 4.3999 g | 20.3591 g | 17.8812 g | | | |
| 240-57930-A-5 | F-VB-UST-P-1 | Moisture | T | 4.3999 g | 20.9501 g | 18.4096 g | | | |
| 240-57930-A-4 | F-VS-UST-P-4 | Moisture | T | 4.3999 g | 26.5689 g | 22.8205 g | | | |
| 240-57930-A-3 | F-VS-UST-P-3 | Moisture | T | 4.3999 g | 32.9572 g | 28.5613 g | | | |
| 240-57930-A-2 | F-VS-UST-P-2 | Moisture | T | 4.3999 g | 32.1604 g | 27.8729 g | | | |
| 240-57930-A-1 | F-VS-UST-P-1 | Moisture | T | 4.3999 g | 25.5114 g | 22.3368 g | | | |

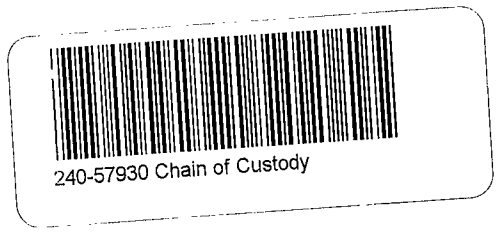
| Batch Notes | |
|---|---|
| Balance ID | B047 No Unit |
| Date and Time Samples in Desiccator | 11/17/15; 05:00 |
| Date and Time Samples out of Desiccator | 11/17/15; 9:35 |
| Date samples were placed in the oven | 11/16/15 |
| Oven Temp when samples are put in oven | 103.9 Degrees C |
| Time samples were place in the oven | 16:30 |
| Date samples were removed from oven | 11/17/15 |
| Oven Temp when samples removed from oven | 103.6 Degrees C |
| Time Samples were removed from oven | 05:00 |
| Oven ID | 002 |
| ID number of the thermometer | Tempguard Box-C #6 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Shipping and Receiving Documents

# CHAIN OF CUSTODY
# AND
# RECEIVING DOCUMENTS





4101 Shuffel Street, N.W. North Canton, OH 44720 **tel** 330.497.9396 **fax** 330.497.0772 **www.testamericainc.com**

3.3/C3.4

# Chain of Custody Record

#48 016105

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING
TestAmerica Laboratories, Inc.
TAL-8210 (0713)

Regulatory Program: ☐ DW ☐ NPDES ☐ RCRA ☐ Other:

| Client Contact | Project Manager: SCOTT NESBIT | Site Contact: MULLIS | Date: 11/13/15 | COC No: |
|---|---|---|---|---|
| Company Name: TETRA TECH | Tel/Fax: 412 921 7134 | Lab Contact: MCFADDEN | Carrier: FedEx | 1 of 1 COCs |
| Address: 20251 Century Blvd. Ste. 200 | Analysis Turnaround Time | | | Sampler: |
| City/State/Zip: Germantown, MD, 20874 | ☐ CALENDAR DAYS ☐ WORKING DAYS | | | For Lab Use Only: |
| Phone: 301-528-3000 | TAT if different from Below | | | Walk-in Client: |
| Fax: 301-528-3000 | ☐ 2 weeks (5 DAYS) | | | Lab Sampling: |
| Project Name: BLOCK F, MRC SOIL REMEDY | ☐ 1 week | | | |
| Site: BLOCK F, UST PIPE INV. | ☐ 2 days | | | Job / SDG No.: |
| PO # 112IC07336 | ☐ 1 day | | | |

| Sample Identification | Sample Date | Sample Time | Sample Type (C=Comp, G=Grab) | Matrix | # of Cont. | Filtered Sample (Y/N) | Perform MS / MSD (Y/N) | VOCs (8260) | TPH GRO/DRO (8015) | SVOCs (8270) | PCBs (8082) | Sample Specific Notes: |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| F-VS-UST-P-1 | 11/13/15 | 1030 | G | SO | 3 | | | K | K | K | K | SAMPLES MAY |
| F-VS-UST-P-2 | | 1035 | | | | | | K | K | K | K | CONTAIN HIGH |
| F-VS-UST-P-3 | | 1040 | | | | | | K | K | K | K | CONTAMINATION |
| F-VS-UST-P-4 | | 1045 | | | | | | K | K | K | K | |
| F-VB-UST-P-1 | | 1000 | | | | | | K | K | K | K | |
| F-VB-UST-P-2 | | 1005 | | | | | | K | K | K | K | |
| F-VB-UST-P-3 | | 1010 | | | | | | K | K | K | K | |
| F-VB-UST-P-4 | | 1015 | | | | | | K | K | K | K | |
| F-VB-UST-P-5 | | 1020 | | | | | | K | K | K | K | |
| F-VB-UST-P-6 | | 1025 | | | | | | K | K | K | K | |
| F-VS-UST-P-DUP | | — | | | | | | K | K | K | K | # DUPLICATE |
| TB-111315 | | — | | AQ | 2 | | | K | | | | |

Page 822 of 823

Preservation Used: 1= Ice, 2= HCl; 3= H2SO4; 4=HNO3; 5=NaOH; 6= Other _____

**Possible Hazard Identification:**
Are any samples from a listed EPA Hazardous Waste? Please List any EPA Waste Codes for the sample in the Comments Section if the lab is to dispose of the sample.

☒ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

**Sample Disposal ( A fee may be assessed if samples are retained longer than 1 month)**

☐ Return to Client ☒ Disposal by Lab ☐ Archive for _____ Months

**Special Instructions/QC Requirements & Comments:** MRC SOIL REMEDY; BLOCK F UST PIPE INVESTIGATION SAMPLES
# VS Samples may contain high contamination, protect instruments

Custody Seals Intact: ☐ Yes ☐ No | Custody Seal No.: | Cooler Temp. (°C): Obs'd: _____ Corr'd: _____ | Therm ID No.: _____

| | Company | Date/Time | | Company | Date/Time |
|---|---|---|---|---|---|
| Relinquished by: | Tetra Tech | 11/13/15 1500 | Received by: | TestAmerica | 11/13/15 1500 |
| Relinquished by: | TT | 11/13/15 1680 | Received by: | | |
| Relinquished by: | | | Received in Laboratory by: Jacquelyn Jenel | TANC | 11/14/15 945 |

**TestAmerica Canton Sample Receipt Form/Narrative** **Login # : 57930**
**Canton Facility**

Client Tetra Tech Site Name _____ Cooler unpacked by: J. Ferrel

Cooler Received on 11/14/15 Opened on 11/14/15

FedEx: 1st Grd (Exp) UPS FAS Stetson Client Drop Off TestAmerica Courier Other _____

**Receipt After-hours**: Drop-off Date/Time _____ Storage Location _____

TestAmerica Cooler # _____ Foam Box (Client Cooler) Box Other _____

Packing material used: (Bubble Wrap) Foam Plastic Bag None Other _____

COOLANT: (Wet Ice) Blue Ice Dry Ice Water None

1. Cooler temperature upon receipt

IR GUN# 53 (CF +**0.1** °C) Observed Cooler Temp. 3.3 °C Corrected Cooler Temp. 3.4 °C

IR GUN# 48 (CF **-0.3** °C) Observed Cooler Temp. _____ °C Corrected Cooler Temp. _____ °C ☐ See Multiple Cooler Form

~~IR GUN# 5 (CF +0.4 °C) Observed Cooler Temp. _____ °C Corrected Cooler Temp. _____ °C~~

IR GUN# 8 (CF **-0.5** °C) Observed Cooler Temp. _____ °C Corrected Cooler Temp. _____ °C

2. Were custody seals on the outside of the cooler(s)? If Yes Quantity 1 (Yes) No
   -Were custody seals on the outside of the cooler(s) signed & dated? (Yes) No NA
   -Were custody seals on the bottle(s) or bottle kits (LLHg/MeHg)? 4CF (Yes) No
3. Shippers' packing slip attached to the cooler(s)? 11/14/15 (Yes) No
4. Did custody papers accompany the sample(s)? (Yes) No
5. Were the custody papers relinquished & signed in the appropriate place? (Yes) No
6. Was/were the person(s) who collected the samples clearly identified on the COC? Yes (No)
7. Did all bottles arrive in good condition (Unbroken)? (Yes) No
8. Could all bottle labels be reconciled with the COC? (Yes) No
9. Were correct bottle(s) used for the test(s) indicated? (Yes) No
10. Sufficient quantity received to perform indicated analyses? (Yes) No
11. Were sample(s) at the correct pH upon receipt? Yes No (NA) pH Strip Lot# **HC554612**
12. Were VOAs on the COC? (Yes) No
13. Were air bubbles >6 mm in any VOA vials? Yes (No) NA
14. Was a trip blank present in the cooler(s)? Trip Blank Lot # NONE (Yes) No

Contacted PM _____ Date _____ by _____ via Verbal Voice Mail Other

Concerning _____

**14. CHAIN OF CUSTODY & SAMPLE DISCREPANCIES** Samples processed by:

**15. SAMPLE CONDITION**

Sample(s) _____ were received after the recommended holding time had expired.

Sample(s) _____ were received in a broken container.

Sample(s) _____ were received with bubble >6 mm in diameter. (Notify PM)

**16. SAMPLE PRESERVATION**

Sample(s) _____ were further preserved in the laboratory.

Time preserved: _____ Preservative(s) added/Lot number(s): _____

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

# ANALYTICAL REPORT

Job Number: 240-59482-1

Job Description: MRC Block F UST Pipe Inv.

For:
Tetra Tech, Inc.
Foster Plaza 7
661 Anderson Drive
Pittsburgh, PA 15220-2745

Attention: Mr. Scott Nesbit

Approved for release.
John McFadden
Project Manager I
1/6/2016 8:34 AM

John McFadden, Project Manager I
4101 Shuffel Street NW, North Canton, OH, 44720
john.mcfadden@testamericainc.com
01/06/2016

cc: Tony Apanavage
Samantha Brenner
Kelly Carper
Tobrena Sedlmyer

The test results in this report meet all NELAP requirements for parameters for which accreditation is required or available. Any exceptions to the NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory. This report is confidential and is intended for the sole use of TestAmerica and its client. All questions regarding this report should be directed to the TestAmerica Project Manager who has signed this report.

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# CASE NARRATIVE

## Client: Tetra Tech, Inc.

## Project: MRC Block F UST Pipe Inv.

## Report Number: 240-59482-1

With the exceptions noted as flags or footnotes, standard analytical protocols were followed in the analysis of the samples and no problems were encountered or anomalies observed. In addition all laboratory quality control samples were within established control limits, with any exceptions noted below. Each sample was analyzed to achieve the lowest possible reporting limit within the constraints of the method. In some cases, due to interference or analytes present at high concentrations, samples were diluted. For diluted samples, the reporting limits are adjusted relative to the dilution required.

TestAmerica Canton attests to the validity of the laboratory data generated by TestAmerica facilities reported herein. All analyses performed by TestAmerica facilities were done using established laboratory SOPs that incorporate QA/QC procedures described in the application methods. TestAmerica's operations groups have reviewed the data for compliance with the laboratory QA/QC plan, and data have been found to be compliant with laboratory protocols unless otherwise noted below.

The test results in this report meet all NELAP requirements for parameters for which accreditation is required or available. Any exceptions to NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory.

Calculations are performed before rounding to avoid round-off errors in calculated results.

All holding times were met and proper preservation noted for the methods performed on these samples, unless otherwise detailed in the individual sections below.

All solid sample results are reported on an "as received" basis unless otherwise indicated by the presence of a % solids value in the method header.

This laboratory report is confidential and is intended for the sole use of TestAmerica and its client.

**RECEIPT**
The samples were received on 12/29/2015 9:30 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 2.5º C.

**VOLATILE ORGANIC COMPOUNDS (GCMS)**
Samples F-VB-UST-P-7 (240-59482-2) and F-VB-UST-P-8 (240-59482-3) were analyzed for volatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8260C. The samples were prepared and analyzed on 12/30/2015.

1,2,3-Trichlorobenzene, 1,2,4-Trichlorobenzene and Naphthalene were detected in method blank MB 240-212606/24 at levels that were above the method detection limit but below the reporting limit. The values should be considered estimates, and have been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**VOLATILE ORGANIC COMPOUNDS (GCMS)**
Sample TB-122815 (240-59482-1) was analyzed for volatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8260C. The samples were analyzed on 01/04/2016.

2-Butanone (MEK) was detected in method blank MB 240-212893/8 at a level that was above the method detection limit but below the reporting limit. The value should be considered an estimate, and has been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

The continuing calibration verification (CCV) for analytical batch 212893 exceeded control criteria for multiple compounds. The samples associated with this CCV were non-detects for the affected analytes. In accordance with the laboratory SOP, a low level CCV at the reporting limit (labeled as an MRL) was analyzed and the affected compounds were detected; therefore the data has been reported. No further corrective action was required. TB-122815 (240-59482-1)

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**SEMIVOLATILE ORGANIC COMPOUNDS (GCMS)**
Samples F-VB-UST-P-7 (240-59482-2) and F-VB-UST-P-8 (240-59482-3) were analyzed for semivolatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8270D. The samples were prepared on 12/30/2015 and analyzed on 12/31/2015.

Bis(2-ethylhexyl) phthalate, Di-n-butyl phthalate and Isophorone were detected in method blank MB 240-212548/7-A at levels that were above the method detection limit but below the reporting limit. The values should be considered estimates, and have been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

The continuing calibration verification (CCV) associated with analytical batch 212723 recovered above the upper control limit for 4-Nitrophenol. The samples associated with this CCV were non-detect for this analyte; therefore, the data have been reported. The following samples are impacted: F-VB-UST-P-7 (240-59482-2), F-VB-UST-P-8 (240-59482-3) and (MB 240-212548/7-A).

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**GASOLINE RANGE ORGANICS**

Samples F-VB-UST-P-7 (240-59482-2) and F-VB-UST-P-8 (240-59482-3) were analyzed for gasoline range organics in accordance with EPA SW-846 Method 8015C - GRO. The samples were prepared and analyzed on 12/30/2015.

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**DIESEL RANGE ORGANICS (DRO)**

Samples F-VB-UST-P-7 (240-59482-2) and F-VB-UST-P-8 (240-59482-3) were analyzed for diesel range organics (DRO) in accordance with EPA SW-846 Method 8015C - DRO. The samples were prepared on 12/30/2015 and analyzed on 12/31/2015.

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**POLYCHLORINATED BIPHENYLS (PCBS)**

Samples F-VB-UST-P-7 (240-59482-2) and F-VB-UST-P-8 (240-59482-3) were analyzed for polychlorinated biphenyls (PCBs) in accordance with EPA SW-846 Method 8082A. The samples were prepared on 12/30/2015 and analyzed on 01/04/2016.

Surrogates are added during the extraction process prior to dilution. When the sample dilution is 5X or greater, surrogate recoveries are diluted out and no corrective action is required. All of the samples in this data set analyzed for PCBs were subjected to the sulfuric acid cleanup procedure before instrumental analysis, per EPA Method 3665A.

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**PERCENT SOLIDS**

Samples F-VB-UST-P-7 (240-59482-2) and F-VB-UST-P-8 (240-59482-3) were analyzed for percent solids in accordance with EPA Method 160.3 MOD. The samples were analyzed on 12/29/2015.

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A4HP9 Analysis Batch Number: 211933

Lab Sample ID: STD4 240-211933/3 IC Client Sample ID:

Date Analyzed: 12/22/15 13:13 Lab File ID: 51222003.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.41 | Poor chromatography | ulmanm | 12/22/15 13:55 |
| Benzoic acid | 7.36 | Poor chromatography | ulmanm | 12/22/15 13:55 |
| 4-Chloroaniline | 7.71 | Poor chromatography | ulmanm | 12/22/15 13:55 |

Lab Sample ID: STD3 240-211933/4 IC Client Sample ID:

Date Analyzed: 12/22/15 13:39 Lab File ID: 51222004.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.42 | Poor chromatography | ulmanm | 12/22/15 14:27 |

Lab Sample ID: STD2 240-211933/5 IC Client Sample ID:

Date Analyzed: 12/22/15 14:04 Lab File ID: 51222005.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 1,4-Dioxane | 3.40 | Poor chromatography | ulmanm | 12/22/15 14:44 |
| Benzoic acid | 7.34 | Poor chromatography | ulmanm | 12/22/15 14:44 |
| 4-Chloroaniline | 7.71 | Poor chromatography | ulmanm | 12/22/15 14:44 |
| Dibenz(a,h)anthracene | 16.91 | Poor chromatography | ulmanm | 12/22/15 14:44 |

Lab Sample ID: STD1 240-211933/6 IC Client Sample ID:

Date Analyzed: 12/22/15 14:30 Lab File ID: 51222006.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Indeno[1,2,3-cd]pyrene | 16.92 | Poor chromatography | ulmanm | 12/22/15 15:05 |
| Dibenz(a,h)anthracene | 16.93 | Poor chromatography | ulmanm | 12/22/15 15:05 |
| Benzo[g,h,i]perylene | 17.59 | Poor chromatography | ulmanm | 12/22/15 15:05 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A4HP9 Analysis Batch Number: 211933

Lab Sample ID: STD9 240-211933/7 IC Client Sample ID:

Date Analyzed: 12/22/15 14:56 Lab File ID: 51222007.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Caprolactam | 8.02 | Poor chromatography | ulmanm | 12/22/15 15:28 |

Lab Sample ID: STD8 240-211933/8 IC Client Sample ID:

Date Analyzed: 12/22/15 15:21 Lab File ID: 51222008.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Caprolactam | 8.02 | Poor chromatography | ulmanm | 12/23/15 07:05 |

Lab Sample ID: STD6 240-211933/10 ICIS Client Sample ID:

Date Analyzed: 12/22/15 16:12 Lab File ID: 51222010.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| 4-Chloroaniline | 7.72 | Poor chromatography | ulmanm | 12/23/15 07:13 |

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A4HP9 Analysis Batch Number: 212723

Lab Sample ID: 240-59482-3 Client Sample ID: F-VB-UST-P-8

Date Analyzed: 12/31/15 12:08 Lab File ID: 51231007.D GC Column: RXI-5SILMS ID: 0.45(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| Benzo[k]fluoranthene | 13.97 | Poor chromatography | ulmanm | 12/31/15 12:48 |
| Dibenz(a,h)anthracene | 16.55 | Poor chromatography | ulmanm | 01/04/16 07:23 |

PCBS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A2HP12 Analysis Batch Number: 207007

Lab Sample ID: STD05 240-207007/7 IC Client Sample ID:

Date Analyzed: 11/16/15 21:09 Lab File ID: P1200042.D GC Column: CLP-1 ID: 0.53(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| PCB-1262 Peak 1 | 7.95 | Baseline Smoothing | hassl | 11/17/15 06:39 |
| PCB-1262 Peak 2 | 8.75 | Baseline Smoothing | hassl | 11/17/15 06:39 |
| PCB-1262 Peak 3 | 9.03 | Baseline Smoothing | hassl | 11/17/15 06:39 |
| PCB-1262 Peak 4 | 9.40 | Baseline Smoothing | hassl | 11/17/15 06:39 |
| PCB-1262 Peak 5 | 9.77 | Baseline Smoothing | hassl | 11/17/15 06:39 |

Lab Sample ID: STD1 240-207007/8 IC Client Sample ID:

Date Analyzed: 11/16/15 21:27 Lab File ID: P1200043.D GC Column: CLP-1 ID: 0.53(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| PCB-1262 Peak 1 | 7.95 | Baseline Smoothing | hassl | 11/17/15 06:40 |
| PCB-1262 Peak 2 | 8.75 | Baseline Smoothing | hassl | 11/17/15 06:40 |
| PCB-1262 Peak 3 | 9.03 | Baseline Smoothing | hassl | 11/17/15 06:40 |
| PCB-1262 Peak 4 | 9.40 | Baseline Smoothing | hassl | 11/17/15 06:40 |
| PCB-1262 Peak 5 | 9.77 | Baseline Smoothing | hassl | 11/17/15 06:40 |

Lab Sample ID: STD02 240-207007/30 IC Client Sample ID:

Date Analyzed: 11/17/15 04:05 Lab File ID: P1200065.D GC Column: CLP-1 ID: 0.53(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| PCB-1016 Peak 1 | 4.61 | Baseline Smoothing | hassl | 11/17/15 06:42 |
| PCB-1016 Peak 2 | 5.11 | Baseline Smoothing | hassl | 11/17/15 06:42 |
| PCB-1016 Peak 3 | 5.74 | Baseline Smoothing | hassl | 11/17/15 06:42 |
| PCB-1016 Peak 4 | 5.91 | Baseline Smoothing | hassl | 11/17/15 06:42 |
| PCB-1016 Peak 5 | 6.45 | Baseline Smoothing | hassl | 11/17/15 06:42 |

PCBS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A2HP12 Analysis Batch Number: 207007

Lab Sample ID: ICV 240-207007/39 Client Sample ID:

Date Analyzed: 11/17/15 06:48 Lab File ID: P1200074.D GC Column: CLP-1 ID: 0.53(mm)

| COMPOUND NAME | RETENTION TIME | MANUAL INTEGRATION | | |
|---|---|---|---|---|
| | | REASON | ANALYST | DATE |
| PCB-1260 Peak 1 | 7.95 | Baseline Smoothing | hassl | 11/17/15 08:38 |
| PCB-1260 Peak 2 | 8.29 | Baseline Smoothing | hassl | 11/17/15 08:38 |
| PCB-1260 Peak 3 | 8.61 | Baseline Smoothing | hassl | 11/17/15 08:38 |
| PCB-1260 Peak 4 | 9.40 | Baseline Smoothing | hassl | 11/17/15 08:38 |
| PCB-1260 Peak 5 | 9.72 | Baseline Smoothing | hassl | 11/17/15 08:38 |

# SAMPLE SUMMARY

Client: Tetra Tech, Inc. Job Number: 240-59482-1

| Lab Sample ID | Client Sample ID | Client Matrix | Date/Time Sampled | Date/Time Received |
|---|---|---|---|---|
| 240-59482-1 | TB-122815 | Water | 12/28/2015 0000 | 12/29/2015 0930 |
| 240-59482-2 | F-VB-UST-P-7 | Solid | 12/28/2015 1000 | 12/29/2015 0930 |
| 240-59482-3 | F-VB-UST-P-8 | Solid | 12/28/2015 1030 | 12/29/2015 0930 |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-59482-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-59482-1** | **TB-122815** | | | | | |
| 2-Butanone (MEK) | | 1.5 | J B | 10 | ug/L | 8260C |
| Acetone | | 1.8 | J | 10 | ug/L | 8260C |
| tert-Butyl alcohol (TBA) | | 5.1 | J | 50 | ug/L | 8260C |
| **240-59482-2** | **F-VB-UST-P-7** | | | | | |
| 1,2,3-Trimethylbenzene | | 0.86 | J | 5.8 | ug/Kg | 8260C |
| 1,2,4-Trimethylbenzene | | 1.9 | J | 5.8 | ug/Kg | 8260C |
| 2-Butanone (MEK) | | 4.8 | J | 23 | ug/Kg | 8260C |
| Acetone | | 36 | | 23 | ug/Kg | 8260C |
| Ethylbenzene | | 0.66 | J | 5.8 | ug/Kg | 8260C |
| Isopropylbenzene | | 0.24 | J F2 | 5.8 | ug/Kg | 8260C |
| m-Xylene & p-Xylene | | 2.9 | J | 12 | ug/Kg | 8260C |
| Naphthalene | | 0.61 | J B | 5.8 | ug/Kg | 8260C |
| o-Xylene | | 1.2 | J | 5.8 | ug/Kg | 8260C |
| Toluene | | 1.6 | J | 5.8 | ug/Kg | 8260C |
| Xylenes, Total | | 4.1 | J | 12 | ug/Kg | 8260C |
| Bis(2-ethylhexyl) phthalate | | 160 | B | 86 | ug/Kg | 8270D |
| Di-n-butyl phthalate | | 30 | J B F1 | 86 | ug/Kg | 8270D |
| Fluoranthene | | 9.3 | | 8.2 | ug/Kg | 8270D |
| Isophorone | | 21 | J B | 61 | ug/Kg | 8270D |
| Naphthalene | | 4.3 | J | 8.2 | ug/Kg | 8270D |
| Phenanthrene | | 9.3 | | 8.2 | ug/Kg | 8270D |
| Pyrene | | 7.9 | J | 8.2 | ug/Kg | 8270D |
| Percent Solids | | 82 | | 0.10 | % | Moisture |
| Percent Moisture | | 18 | | 0.10 | % | Moisture |

# EXECUTIVE SUMMARY - Detections

Client: Tetra Tech, Inc. Job Number: 240-59482-1

| Lab Sample ID / Analyte | Client Sample ID | Result | Qualifier | Reporting Limit | Units | Method |
|---|---|---|---|---|---|---|
| **240-59482-3** | **F-VB-UST-P-8** | | | | | |
| 2-Butanone (MEK) | | 1.9 | J | 22 | ug/Kg | 8260C |
| p-Isopropyltoluene | | 0.57 | J | 5.5 | ug/Kg | 8260C |
| Acetone | | 9.4 | J | 22 | ug/Kg | 8260C |
| Isopropylbenzene | | 2.0 | J | 5.5 | ug/Kg | 8260C |
| Naphthalene | | 0.88 | J B | 5.5 | ug/Kg | 8260C |
| n-Butylbenzene | | 0.83 | J | 5.5 | ug/Kg | 8260C |
| N-Propylbenzene | | 1.9 | J | 5.5 | ug/Kg | 8260C |
| 2-Methylnaphthalene | | 53 | | 7.9 | ug/Kg | 8270D |
| Acenaphthene | | 63 | | 7.9 | ug/Kg | 8270D |
| Anthracene | | 97 | | 7.9 | ug/Kg | 8270D |
| Benzo[a]anthracene | | 140 | | 7.9 | ug/Kg | 8270D |
| Benzo[a]pyrene | | 100 | | 7.9 | ug/Kg | 8270D |
| Benzo[b]fluoranthene | | 180 | | 7.9 | ug/Kg | 8270D |
| Benzo[g,h,i]perylene | | 51 | | 7.9 | ug/Kg | 8270D |
| Benzo[k]fluoranthene | | 47 | | 7.9 | ug/Kg | 8270D |
| Bis(2-ethylhexyl) phthalate | | 82 | B | 82 | ug/Kg | 8270D |
| Chrysene | | 94 | | 7.9 | ug/Kg | 8270D |
| Dibenz(a,h)anthracene | | 16 | | 7.9 | ug/Kg | 8270D |
| Dibenzofuran | | 51 | J | 59 | ug/Kg | 8270D |
| Di-n-butyl phthalate | | 23 | J B | 82 | ug/Kg | 8270D |
| Fluoranthene | | 490 | | 7.9 | ug/Kg | 8270D |
| Fluorene | | 88 | | 7.9 | ug/Kg | 8270D |
| Indeno[1,2,3-cd]pyrene | | 45 | | 7.9 | ug/Kg | 8270D |
| Isophorone | | 20 | J B | 59 | ug/Kg | 8270D |
| Naphthalene | | 14 | | 7.9 | ug/Kg | 8270D |
| Phenanthrene | | 570 | | 7.9 | ug/Kg | 8270D |
| Pyrene | | 370 | | 7.9 | ug/Kg | 8270D |
| Total Petroleum Hydrocarbons (C10-C32) | | 42 | | 20 | mg/Kg | 8015C |
| Percent Solids | | 84 | | 0.10 | % | Moisture |
| Percent Moisture | | 16 | | 0.10 | % | Moisture |

# METHOD SUMMARY

Client: Tetra Tech, Inc. Job Number: 240-59482-1

| Description | Lab Location | Method | Preparation Method |
|---|---|---|---|
| **Matrix: Solid** | | | |
| Volatile Organic Compounds by GC/MS | TAL CAN | SW846 8260C | |
| Purge and Trap | TAL CAN | | SW846 5030C |
| Semivolatile Organic Compounds (GC/MS) | TAL CAN | SW846 8270D | |
| Soxhlet Extraction | TAL CAN | | SW846 3540C |
| Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics) | TAL CAN | SW846 8015C | |
| Purge and Trap | TAL CAN | | SW846 5030A |
| Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics) | TAL CAN | SW846 8015C | |
| Soxhlet Extraction | TAL CAN | | SW846 3540C |
| Polychlorinated Biphenyls (PCBs) by Gas Chromatography | TAL CAN | SW846 8082A | |
| Soxhlet Extraction | TAL CAN | | SW846 3540C |
| Percent Moisture | TAL CAN | EPA Moisture | |
| **Matrix: Water** | | | |
| Volatile Organic Compounds by GC/MS | TAL CAN | SW846 8260C | |
| Purge and Trap | TAL CAN | | SW846 5030C |

### Lab References:

TAL CAN = TestAmerica Canton

### Method References:

EPA = US Environmental Protection Agency

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

# METHOD / ANALYST SUMMARY

Client: Tetra Tech, Inc. Job Number: 240-59482-1

| Method | Analyst | Analyst ID |
|---|---|---|
| SW846 8260C | Evans, Laura | LEE |
| SW846 8260C | Macenczak, Steven | SAM |
| SW846 8270D | Ulman, Mark | MRU |
| SW846 8015C | Reed, Rachel T | RTR |
| SW846 8015C | Bosworth, Heather M | HMB |
| SW846 8082A | Bosworth, Heather M | HMB |
| EPA Moisture | Woodward, Bruce | BLW |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID: TB-122815**

Lab Sample ID: 240-59482-1 Date Sampled: 12/28/2015 0000
Client Matrix: Water Date Received: 12/29/2015 0930

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Prep Method: | 5030C | Prep Batch: | N/A | Lab File ID: | UXC5597.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1207 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1207 | | | | |

| Analyte | Result (ug/L) | Qualifier | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.28 | U | 0.28 | 1.0 |
| 1,1,1-Trichloroethane | 0.44 | U | 0.44 | 1.0 |
| 1,1,2,2-Tetrachloroethane | 0.22 | U | 0.22 | 1.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloroethane | 0.30 | U | 0.30 | 1.0 |
| 1,1-Dichloroethene | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloropropene | 0.42 | U | 0.42 | 1.0 |
| 1,2,3-Trichlorobenzene | 0.37 | U | 0.37 | 1.0 |
| 1,2,3-Trichloropropane | 0.44 | U | 0.44 | 1.0 |
| 1,2,3-Trimethylbenzene | 0.47 | U | 0.47 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.32 | U | 0.32 | 1.0 |
| 1,2,4-Trimethylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromo-3-Chloropropane | 0.82 | U | 0.82 | 2.0 |
| 1,2-Dichlorobenzene | 0.25 | U | 0.25 | 1.0 |
| 1,2-Dichloroethane | 0.23 | U | 0.23 | 1.0 |
| 1,2-Dichloropropane | 0.25 | U | 0.25 | 1.0 |
| 1,3-Dichlorobenzene | 0.19 | U | 0.19 | 1.0 |
| 1,3-Dichloropropane | 0.19 | U | 0.19 | 1.0 |
| 1,4-Dichlorobenzene | 0.27 | U | 0.27 | 1.0 |
| 2,2-Dichloropropane | 0.26 | U | 0.26 | 1.0 |
| 2-Butanone (MEK) | 1.5 | J B | 0.53 | 10 |
| 2-Chloroethyl vinyl ether | 0.59 | U | 0.59 | 10 |
| 2-Chlorotoluene | 0.40 | U | 0.40 | 1.0 |
| 2-Hexanone | 0.48 | U | 0.48 | 10 |
| 4-Chlorotoluene | 0.29 | U | 0.29 | 1.0 |
| p-Isopropyltoluene | 0.43 | U | 0.43 | 1.0 |
| 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 0.99 | 10 |
| Acetone | 1.8 | J | 0.94 | 10 |
| Benzene | 0.35 | U | 0.35 | 1.0 |
| Bromobenzene | 0.35 | U | 0.35 | 1.0 |
| Bromoform | 0.56 | U | 0.56 | 1.0 |
| Bromomethane | 0.44 | U | 0.44 | 1.0 |
| Carbon disulfide | 0.38 | U | 0.38 | 1.0 |
| Carbon tetrachloride | 0.43 | U | 0.43 | 1.0 |
| Chlorobenzene | 0.25 | U | 0.25 | 1.0 |
| Dibromochloromethane | 0.43 | U | 0.43 | 1.0 |
| Chloroethane | 0.32 | U | 0.32 | 1.0 |
| Chloroform | 0.25 | U | 0.25 | 1.0 |
| Chloromethane | 0.44 | U | 0.44 | 1.0 |
| cis-1,2-Dichloroethene | 0.26 | U | 0.26 | 1.0 |
| cis-1,3-Dichloropropene | 0.46 | U | 0.46 | 1.0 |
| Dibromomethane | 0.42 | U | 0.42 | 1.0 |
| Bromochloromethane | 0.50 | U | 0.50 | 1.0 |
| Dichlorodifluoromethane | 0.32 | U | 0.32 | 1.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 1.0 |
| tert-Butyl alcohol (TBA) | 5.1 | J | 4.9 | 50 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **TB-122815**

Lab Sample ID: 240-59482-1 Date Sampled: 12/28/2015 0000
Client Matrix: Water Date Received: 12/29/2015 0930

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Prep Method: | 5030C | Prep Batch: | N/A | Lab File ID: | UXC5597.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1207 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1207 | | | | |

| Analyte | Result (ug/L) | Qualifier | MDL | RL |
|---|---|---|---|---|
| Ethylbenzene | 0.25 | U | 0.25 | 1.0 |
| Hexachlorobutadiene | 0.35 | U | 0.35 | 1.0 |
| Diisopropyl ether | 0.50 | U | 0.50 | 10 |
| Isopropylbenzene | 0.35 | U | 0.35 | 1.0 |
| Methyl tert-butyl ether | 0.20 | U | 0.20 | 1.0 |
| Methylene Chloride | 0.33 | U | 0.33 | 1.0 |
| m-Xylene & p-Xylene | 0.24 | U | 0.24 | 2.0 |
| Naphthalene | 0.45 | U | 0.45 | 1.0 |
| n-Butylbenzene | 0.31 | U | 0.31 | 1.0 |
| N-Propylbenzene | 0.40 | U | 0.40 | 1.0 |
| o-Xylene | 0.25 | U | 0.25 | 1.0 |
| sec-Butylbenzene | 0.48 | U | 0.48 | 1.0 |
| Styrene | 0.45 | U | 0.45 | 1.0 |
| Tert-amyl methyl ether | 0.30 | U | 0.30 | 5.0 |
| Ethyl t-butyl ether | 0.23 | U | 0.23 | 5.0 |
| Tetrachloroethene | 0.31 | U | 0.31 | 1.0 |
| Toluene | 0.23 | U | 0.23 | 1.0 |
| trans-1,2-Dichloroethene | 0.30 | U | 0.30 | 1.0 |
| trans-1,3-Dichloropropene | 0.56 | U | 0.56 | 1.0 |
| Trichloroethene | 0.22 | U | 0.22 | 1.0 |
| Trichlorofluoromethane | 0.49 | U | 0.49 | 1.0 |
| Vinyl acetate | 0.41 | U | 0.41 | 2.0 |
| Vinyl chloride | 0.29 | U | 0.29 | 1.0 |
| Xylenes, Total | 0.52 | U | 0.52 | 2.0 |
| tert-Butylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromoethane | 0.32 | U | 0.32 | 1.0 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 102 | | 80 - 120 |
| Dibromofluoromethane (Surr) | 93 | | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 112 | | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 99 | | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-7**

Lab Sample ID: 240-59482-2 Date Sampled: 12/28/2015 1000
Client Matrix: Solid % Moisture: 18.4 Date Received: 12/29/2015 0930

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-212657 | Lab File ID: | 147326.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.26 g |
| Analysis Date: | 12/30/2015 2148 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.58 | U | 0.58 | 5.8 |
| 1,1,1-Trichloroethane | | 0.76 | U | 0.76 | 5.8 |
| 1,1,2,2-Tetrachloroethane | | 0.36 | U | 0.36 | 5.8 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.49 | U | 0.49 | 5.8 |
| 1,1-Dichloroethane | | 0.40 | U | 0.40 | 5.8 |
| 1,1-Dichloroethene | | 0.93 | U | 0.93 | 5.8 |
| 1,1-Dichloropropene | | 0.43 | U | 0.43 | 5.8 |
| 1,2,3-Trichlorobenzene | | 0.58 | U | 0.58 | 5.8 |
| 1,2,3-Trichloropropane | | 0.91 | U | 0.91 | 5.8 |
| 1,2,3-Trimethylbenzene | | 0.86 | J | 0.54 | 5.8 |
| 1,2,4-Trichlorobenzene | | 0.44 | U | 0.44 | 5.8 |
| 1,2,4-Trimethylbenzene | | 1.9 | J | 0.68 | 5.8 |
| 1,2-Dibromo-3-Chloropropane | | 3.2 | U | 3.2 | 12 |
| 1,2-Dichlorobenzene | | 0.40 | U | 0.40 | 5.8 |
| 1,2-Dichloroethane | | 0.54 | U | 0.54 | 5.8 |
| 1,2-Dichloropropane | | 0.16 | U | 0.16 | 5.8 |
| 1,3-Dichlorobenzene | | 0.80 | U F2 | 0.80 | 5.8 |
| 1,3-Dichloropropane | | 0.47 | U | 0.47 | 5.8 |
| 1,4-Dichlorobenzene | | 0.33 | U F2 | 0.33 | 5.8 |
| 2,2-Dichloropropane | | 0.51 | U | 0.51 | 5.8 |
| 2-Butanone (MEK) | | 4.8 | J | 1.3 | 23 |
| 2-Chloroethyl vinyl ether | | 0.75 | U | 0.75 | 58 |
| 2-Chlorotoluene | | 0.35 | U F2 | 0.35 | 5.8 |
| 2-Hexanone | | 0.79 | U | 0.79 | 23 |
| 4-Chlorotoluene | | 0.30 | U F2 | 0.30 | 5.8 |
| p-Isopropyltoluene | | 0.48 | U F2 | 0.48 | 5.8 |
| 4-Methyl-2-pentanone (MIBK) | | 1.4 | U | 1.4 | 23 |
| Acetone | | 36 | | 6.6 | 23 |
| Benzene | | 1.0 | U | 1.0 | 5.8 |
| Bromobenzene | | 0.38 | U | 0.38 | 5.8 |
| Bromoform | | 0.28 | U | 0.28 | 5.8 |
| Bromomethane | | 0.49 | U | 0.49 | 5.8 |
| Carbon disulfide | | 0.69 | U | 0.69 | 5.8 |
| Carbon tetrachloride | | 0.76 | U | 0.76 | 5.8 |
| Chlorobenzene | | 0.62 | U | 0.62 | 5.8 |
| Dibromochloromethane | | 0.43 | U | 0.43 | 5.8 |
| Chloroethane | | 0.51 | U | 0.51 | 5.8 |
| Chloroform | | 0.42 | U | 0.42 | 5.8 |
| Chloromethane | | 0.92 | U | 0.92 | 5.8 |
| cis-1,2-Dichloroethene | | 0.33 | U | 0.33 | 5.8 |
| cis-1,3-Dichloropropene | | 0.83 | U | 0.83 | 5.8 |
| Dibromomethane | | 0.49 | U | 0.49 | 5.8 |
| Bromochloromethane | | 0.38 | U | 0.38 | 5.8 |
| Dichlorodifluoromethane | | 0.33 | U | 0.33 | 5.8 |
| Bromodichloromethane | | 0.34 | U | 0.34 | 5.8 |
| tert-Butyl alcohol (TBA) | | 6.7 | U | 6.7 | 230 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-7**

Lab Sample ID: 240-59482-2 Date Sampled: 12/28/2015 1000
Client Matrix: Solid % Moisture: 18.4 Date Received: 12/29/2015 0930

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-212657 | Lab File ID: | 147326.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.26 g |
| Analysis Date: | 12/30/2015 2148 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.66 | J | 0.37 | 5.8 |
| Hexachlorobutadiene | | 0.47 | U F2 | 0.47 | 5.8 |
| Diisopropyl ether | | 0.69 | U | 0.69 | 12 |
| Isopropylbenzene | | 0.24 | J F2 | 0.23 | 5.8 |
| Methyl tert-butyl ether | | 0.50 | U | 0.50 | 5.8 |
| Methylene Chloride | | 0.86 | U | 0.86 | 5.8 |
| m-Xylene & p-Xylene | | 2.9 | J | 0.49 | 12 |
| Naphthalene | | 0.61 | J B | 0.41 | 5.8 |
| n-Butylbenzene | | 0.69 | U F2 | 0.69 | 5.8 |
| N-Propylbenzene | | 0.70 | U F2 | 0.70 | 5.8 |
| o-Xylene | | 1.2 | J | 0.66 | 5.8 |
| sec-Butylbenzene | | 1.1 | U F2 | 1.1 | 5.8 |
| Styrene | | 0.47 | U | 0.47 | 5.8 |
| Tert-amyl methyl ether | | 0.50 | U | 0.50 | 5.8 |
| Ethyl t-butyl ether | | 0.48 | U | 0.48 | 5.8 |
| Tetrachloroethene | | 0.93 | U | 0.93 | 5.8 |
| Toluene | | 1.6 | J | 0.31 | 5.8 |
| trans-1,2-Dichloroethene | | 0.37 | U | 0.37 | 5.8 |
| trans-1,3-Dichloropropene | | 0.34 | U | 0.34 | 5.8 |
| Trichloroethene | | 0.44 | U | 0.44 | 5.8 |
| Trichlorofluoromethane | | 0.34 | U | 0.34 | 5.8 |
| Vinyl acetate | | 0.77 | U F1 | 0.77 | 12 |
| Vinyl chloride | | 0.35 | U | 0.35 | 5.8 |
| Xylenes, Total | | 4.1 | J | 0.63 | 12 |
| tert-Butylbenzene | | 0.78 | U F2 | 0.78 | 5.8 |
| 1,2-Dibromoethane | | 0.62 | U | 0.62 | 5.8 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 91 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 93 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 88 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 100 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-8**

Lab Sample ID: 240-59482-3 Date Sampled: 12/28/2015 1030
Client Matrix: Solid % Moisture: 16.1 Date Received: 12/29/2015 0930

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-212657 | Lab File ID: | 147330.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.43 g |
| Analysis Date: | 12/30/2015 2314 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | | 0.55 | U | 0.55 | 5.5 |
| 1,1,1-Trichloroethane | | 0.71 | U | 0.71 | 5.5 |
| 1,1,2,2-Tetrachloroethane | | 0.34 | U | 0.34 | 5.5 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | | 0.46 | U | 0.46 | 5.5 |
| 1,1-Dichloroethane | | 0.37 | U | 0.37 | 5.5 |
| 1,1-Dichloroethene | | 0.88 | U | 0.88 | 5.5 |
| 1,1-Dichloropropene | | 0.41 | U | 0.41 | 5.5 |
| 1,2,3-Trichlorobenzene | | 0.55 | U | 0.55 | 5.5 |
| 1,2,3-Trichloropropane | | 0.86 | U | 0.86 | 5.5 |
| 1,2,3-Trimethylbenzene | | 0.50 | U | 0.50 | 5.5 |
| 1,2,4-Trichlorobenzene | | 0.42 | U | 0.42 | 5.5 |
| 1,2,4-Trimethylbenzene | | 0.64 | U | 0.64 | 5.5 |
| 1,2-Dibromo-3-Chloropropane | | 3.0 | U | 3.0 | 11 |
| 1,2-Dichlorobenzene | | 0.37 | U | 0.37 | 5.5 |
| 1,2-Dichloroethane | | 0.50 | U | 0.50 | 5.5 |
| 1,2-Dichloropropane | | 0.15 | U | 0.15 | 5.5 |
| 1,3-Dichlorobenzene | | 0.76 | U | 0.76 | 5.5 |
| 1,3-Dichloropropane | | 0.44 | U | 0.44 | 5.5 |
| 1,4-Dichlorobenzene | | 0.31 | U | 0.31 | 5.5 |
| 2,2-Dichloropropane | | 0.48 | U | 0.48 | 5.5 |
| 2-Butanone (MEK) | | 1.9 | J | 1.2 | 22 |
| 2-Chloroethyl vinyl ether | | 0.70 | U | 0.70 | 55 |
| 2-Chlorotoluene | | 0.33 | U | 0.33 | 5.5 |
| 2-Hexanone | | 0.75 | U | 0.75 | 22 |
| 4-Chlorotoluene | | 0.29 | U | 0.29 | 5.5 |
| p-Isopropyltoluene | | 0.57 | J | 0.45 | 5.5 |
| 4-Methyl-2-pentanone (MIBK) | | 1.3 | U | 1.3 | 22 |
| Acetone | | 9.4 | J | 6.3 | 22 |
| Benzene | | 0.99 | U | 0.99 | 5.5 |
| Bromobenzene | | 0.36 | U | 0.36 | 5.5 |
| Bromoform | | 0.26 | U | 0.26 | 5.5 |
| Bromomethane | | 0.46 | U | 0.46 | 5.5 |
| Carbon disulfide | | 0.65 | U | 0.65 | 5.5 |
| Carbon tetrachloride | | 0.71 | U | 0.71 | 5.5 |
| Chlorobenzene | | 0.58 | U | 0.58 | 5.5 |
| Dibromochloromethane | | 0.41 | U | 0.41 | 5.5 |
| Chloroethane | | 0.48 | U | 0.48 | 5.5 |
| Chloroform | | 0.40 | U | 0.40 | 5.5 |
| Chloromethane | | 0.87 | U | 0.87 | 5.5 |
| cis-1,2-Dichloroethene | | 0.31 | U | 0.31 | 5.5 |
| cis-1,3-Dichloropropene | | 0.78 | U | 0.78 | 5.5 |
| Dibromomethane | | 0.46 | U | 0.46 | 5.5 |
| Bromochloromethane | | 0.36 | U | 0.36 | 5.5 |
| Dichlorodifluoromethane | | 0.31 | U | 0.31 | 5.5 |
| Bromodichloromethane | | 0.32 | U | 0.32 | 5.5 |
| tert-Butyl alcohol (TBA) | | 6.3 | U | 6.3 | 220 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-8**

Lab Sample ID: 240-59482-3 Date Sampled: 12/28/2015 1030
Client Matrix: Solid % Moisture: 16.1 Date Received: 12/29/2015 0930

**8260C Volatile Organic Compounds by GC/MS**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8260C | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Prep Method: | 5030C | Prep Batch: | 240-212657 | Lab File ID: | 147330.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 5.43 g |
| Analysis Date: | 12/30/2015 2314 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Ethylbenzene | | 0.35 | U | 0.35 | 5.5 |
| Hexachlorobutadiene | | 0.44 | U | 0.44 | 5.5 |
| Diisopropyl ether | | 0.65 | U | 0.65 | 11 |
| Isopropylbenzene | | 2.0 | J | 0.22 | 5.5 |
| Methyl tert-butyl ether | | 0.47 | U | 0.47 | 5.5 |
| Methylene Chloride | | 0.81 | U | 0.81 | 5.5 |
| m-Xylene & p-Xylene | | 0.46 | U | 0.46 | 11 |
| Naphthalene | | 0.88 | J B | 0.38 | 5.5 |
| n-Butylbenzene | | 0.83 | J | 0.65 | 5.5 |
| N-Propylbenzene | | 1.9 | J | 0.66 | 5.5 |
| o-Xylene | | 0.63 | U | 0.63 | 5.5 |
| sec-Butylbenzene | | 1.0 | U | 1.0 | 5.5 |
| Styrene | | 0.44 | U | 0.44 | 5.5 |
| Tert-amyl methyl ether | | 0.47 | U | 0.47 | 5.5 |
| Ethyl t-butyl ether | | 0.45 | U | 0.45 | 5.5 |
| Tetrachloroethene | | 0.88 | U | 0.88 | 5.5 |
| Toluene | | 0.30 | U | 0.30 | 5.5 |
| trans-1,2-Dichloroethene | | 0.35 | U | 0.35 | 5.5 |
| trans-1,3-Dichloropropene | | 0.32 | U | 0.32 | 5.5 |
| Trichloroethene | | 0.42 | U | 0.42 | 5.5 |
| Trichlorofluoromethane | | 0.32 | U | 0.32 | 5.5 |
| Vinyl acetate | | 0.72 | U | 0.72 | 11 |
| Vinyl chloride | | 0.33 | U | 0.33 | 5.5 |
| Xylenes, Total | | 0.59 | U | 0.59 | 11 |
| tert-Butylbenzene | | 0.74 | U | 0.74 | 5.5 |
| 1,2-Dibromoethane | | 0.58 | U | 0.58 | 5.5 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 91 | | 67 - 125 |
| Dibromofluoromethane (Surr) | 90 | | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 87 | | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 97 | | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-7**

Lab Sample ID: 240-59482-2 Date Sampled: 12/28/2015 1000
Client Matrix: Solid % Moisture: 18.4 Date Received: 12/29/2015 0930

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-212548 | Lab File ID: | 51231008.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.06 g |
| Analysis Date: | 12/31/2015 1233 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.3 | U | 4.3 | 61 |
| bis (2-chloroisopropyl) ether | | 12 | U | 12 | 120 |
| 2,4,5-Trichlorophenol | | 31 | U | 31 | 180 |
| 2,4,6-Trichlorophenol | | 11 | U | 11 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 26 | U | 26 | 400 |
| 2,4-Dinitrotoluene | | 21 | U | 21 | 240 |
| 2,6-Dinitrotoluene | | 26 | U | 26 | 240 |
| 2-Chloronaphthalene | | 0.55 | U | 0.55 | 61 |
| 2-Chlorophenol | | 10 | U | 10 | 61 |
| 2-Methylnaphthalene | | 0.61 | U | 0.61 | 8.2 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 10 | U | 10 | 61 |
| 3,3'-Dichlorobenzidine | | 22 | U | 22 | 120 |
| 3-Nitroaniline | | 20 | U | 20 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Chloro-3-methylphenol | | 26 | U | 26 | 180 |
| 4-Chloroaniline | | 21 | U | 21 | 180 |
| 4-Chlorophenyl phenyl ether | | 16 | U | 16 | 61 |
| 4-Nitroaniline | | 32 | U | 32 | 240 |
| 4-Nitrophenol | | 21 | U F1 | 21 | 400 |
| Acenaphthene | | 0.93 | U | 0.93 | 8.2 |
| Acenaphthylene | | 0.43 | U | 0.43 | 8.2 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 0.95 | U | 0.95 | 8.2 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 15 | U | 15 | 120 |
| Benzo[a]anthracene | | 0.77 | U | 0.77 | 8.2 |
| Benzo[a]pyrene | | 0.78 | U | 0.78 | 8.2 |
| Benzo[b]fluoranthene | | 0.72 | U | 0.72 | 8.2 |
| Benzo[g,h,i]perylene | | 0.43 | U | 0.43 | 8.2 |
| Benzo[k]fluoranthene | | 0.83 | U | 0.83 | 8.2 |
| Bis(2-chloroethoxy)methane | | 27 | U | 27 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 160 | B | 23 | 86 |
| Butyl benzyl phthalate | | 12 | U | 12 | 86 |
| Caprolactam | | 45 | U | 45 | 400 |
| Carbazole | | 33 | U | 33 | 61 |
| Chrysene | | 1.3 | U | 1.3 | 8.2 |
| Dibenz(a,h)anthracene | | 0.81 | U | 0.81 | 8.2 |
| Dibenzofuran | | 0.81 | U | 0.81 | 61 |
| Diethyl phthalate | | 20 | U | 20 | 86 |
| Dimethyl phthalate | | 21 | U | 21 | 86 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID: F-VB-UST-P-7**

Lab Sample ID: 240-59482-2 Date Sampled: 12/28/2015 1000
Client Matrix: Solid % Moisture: 18.4 Date Received: 12/29/2015 0930

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-212548 | Lab File ID: | 51231008.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.06 g |
| Analysis Date: | 12/31/2015 1233 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 30 | J B F1 | 18 | 86 |
| Di-n-octyl phthalate | | 9.7 | U | 9.7 | 86 |
| Fluoranthene | | 9.3 | | 0.67 | 8.2 |
| Fluorene | | 0.65 | U | 0.65 | 8.2 |
| Hexachlorobenzene | | 2.6 | U | 2.6 | 8.2 |
| Hexachlorobutadiene | | 6.9 | U | 6.9 | 61 |
| Hexachlorocyclopentadiene | | 9.9 | U | 9.9 | 400 |
| Hexachloroethane | | 11 | U | 11 | 61 |
| Indeno[1,2,3-cd]pyrene | | 0.43 | U | 0.43 | 8.2 |
| Isophorone | | 21 | J B | 16 | 61 |
| N-Nitrosodi-n-propylamine | | 7.7 | U | 7.7 | 61 |
| N-Nitrosodiphenylamine | | 26 | U | 26 | 61 |
| Naphthalene | | 4.3 | J | 1.0 | 8.2 |
| Nitrobenzene | | 2.7 | U | 2.7 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 9.3 | | 0.89 | 8.2 |
| Phenol | | 8.9 | U | 8.9 | 61 |
| Pyrene | | 7.9 | J | 0.54 | 8.2 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 490 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 87 | | 36 - 110 |
| Phenol-d5 (Surr) | 83 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 80 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 81 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 77 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 88 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-8**

Lab Sample ID: 240-59482-3 Date Sampled: 12/28/2015 1030
Client Matrix: Solid % Moisture: 16.1 Date Received: 12/29/2015 0930

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-212548 | Lab File ID: | 51231007.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.35 g |
| Analysis Date: | 12/31/2015 1208 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | | 4.1 | U | 4.1 | 59 |
| bis (2-chloroisopropyl) ether | | 11 | U | 11 | 120 |
| 2,4,5-Trichlorophenol | | 29 | U | 29 | 180 |
| 2,4,6-Trichlorophenol | | 10 | U | 10 | 180 |
| 2,4-Dichlorophenol | | 24 | U | 24 | 180 |
| 2,4-Dimethylphenol | | 24 | U | 24 | 180 |
| 2,4-Dinitrophenol | | 25 | U | 25 | 390 |
| 2,4-Dinitrotoluene | | 20 | U | 20 | 240 |
| 2,6-Dinitrotoluene | | 25 | U | 25 | 240 |
| 2-Chloronaphthalene | | 0.53 | U | 0.53 | 59 |
| 2-Chlorophenol | | 9.7 | U | 9.7 | 59 |
| 2-Methylnaphthalene | | 53 | | 0.59 | 7.9 |
| 2-Methylphenol | | 13 | U | 13 | 240 |
| 2-Nitroaniline | | 11 | U | 11 | 240 |
| 2-Nitrophenol | | 9.8 | U | 9.8 | 59 |
| 3,3'-Dichlorobenzidine | | 21 | U | 21 | 120 |
| 3-Nitroaniline | | 19 | U | 19 | 240 |
| 4,6-Dinitro-2-methylphenol | | 11 | U | 11 | 180 |
| 4-Bromophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Chloro-3-methylphenol | | 25 | U | 25 | 180 |
| 4-Chloroaniline | | 20 | U | 20 | 180 |
| 4-Chlorophenyl phenyl ether | | 15 | U | 15 | 59 |
| 4-Nitroaniline | | 31 | U | 31 | 240 |
| 4-Nitrophenol | | 20 | U | 20 | 390 |
| Acenaphthene | | 63 | | 0.90 | 7.9 |
| Acenaphthylene | | 0.41 | U | 0.41 | 7.9 |
| Acetophenone | | 11 | U | 11 | 120 |
| Anthracene | | 97 | | 0.92 | 7.9 |
| Atrazine | | 11 | U | 11 | 240 |
| Benzaldehyde | | 14 | U | 14 | 120 |
| Benzo[a]anthracene | | 140 | | 0.74 | 7.9 |
| Benzo[a]pyrene | | 100 | | 0.75 | 7.9 |
| Benzo[b]fluoranthene | | 180 | | 0.70 | 7.9 |
| Benzo[g,h,i]perylene | | 51 | | 0.41 | 7.9 |
| Benzo[k]fluoranthene | | 47 | | 0.80 | 7.9 |
| Bis(2-chloroethoxy)methane | | 26 | U | 26 | 120 |
| Bis(2-chloroethyl)ether | | 2.4 | U | 2.4 | 120 |
| Bis(2-ethylhexyl) phthalate | | 82 | B | 22 | 82 |
| Butyl benzyl phthalate | | 12 | U | 12 | 82 |
| Caprolactam | | 44 | U | 44 | 390 |
| Carbazole | | 32 | U | 32 | 59 |
| Chrysene | | 94 | | 1.3 | 7.9 |
| Dibenz(a,h)anthracene | | 16 | | 0.78 | 7.9 |
| Dibenzofuran | | 51 | J | 0.78 | 59 |
| Diethyl phthalate | | 19 | U | 19 | 82 |
| Dimethyl phthalate | | 20 | U | 20 | 82 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID: F-VB-UST-P-8**

Lab Sample ID: 240-59482-3 Date Sampled: 12/28/2015 1030
Client Matrix: Solid % Moisture: 16.1 Date Received: 12/29/2015 0930

**8270D Semivolatile Organic Compounds (GC/MS)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8270D | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Prep Method: | 3540C | Prep Batch: | 240-212548 | Lab File ID: | 51231007.D |
| Dilution: | 1.0 | | | Initial Weight/Volume: | 30.35 g |
| Analysis Date: | 12/31/2015 1208 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | | 23 | J B | 18 | 82 |
| Di-n-octyl phthalate | | 9.3 | U | 9.3 | 82 |
| Fluoranthene | | 490 | | 0.65 | 7.9 |
| Fluorene | | 88 | | 0.62 | 7.9 |
| Hexachlorobenzene | | 2.5 | U | 2.5 | 7.9 |
| Hexachlorobutadiene | | 6.6 | U | 6.6 | 59 |
| Hexachlorocyclopentadiene | | 9.5 | U | 9.5 | 390 |
| Hexachloroethane | | 11 | U | 11 | 59 |
| Indeno[1,2,3-cd]pyrene | | 45 | | 0.41 | 7.9 |
| Isophorone | | 20 | J B | 15 | 59 |
| N-Nitrosodi-n-propylamine | | 7.4 | U | 7.4 | 59 |
| N-Nitrosodiphenylamine | | 25 | U | 25 | 59 |
| Naphthalene | | 14 | | 0.97 | 7.9 |
| Nitrobenzene | | 2.6 | U | 2.6 | 120 |
| Pentachlorophenol | | 11 | U | 11 | 180 |
| Phenanthrene | | 570 | | 0.86 | 7.9 |
| Phenol | | 8.6 | U | 8.6 | 59 |
| Pyrene | | 370 | | 0.52 | 7.9 |
| 3 & 4 Methylphenol | | 24 | U | 24 | 470 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 87 | | 36 - 110 |
| Phenol-d5 (Surr) | 85 | | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 81 | | 20 - 110 |
| 2-Fluorophenol (Surr) | 83 | | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 77 | | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 101 | | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-7**

Lab Sample ID: 240-59482-2 Date Sampled: 12/28/2015 1000
Client Matrix: Solid % Moisture: 18.4 Date Received: 12/29/2015 0930

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-212533 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-212559 | Initial Weight/Volume: | 5.05 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 12/30/2015 1023 | | | Injection Volume: | 5 mL |
| Prep Date: | 12/30/2015 0739 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 56 | U | 56 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 70 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID: F-VB-UST-P-8**

Lab Sample ID: 240-59482-3 Date Sampled: 12/28/2015 1030
Client Matrix: Solid % Moisture: 16.1 Date Received: 12/29/2015 0930

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-212533 | Instrument ID: | YPID |
| Prep Method: | 5030A | Prep Batch: | 240-212559 | Initial Weight/Volume: | 5.00 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 12/30/2015 1204 | | | Injection Volume: | 5 mL |
| Prep Date: | 12/30/2015 0739 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | | 55 | U | 55 | 120 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 67 | | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-7**

Lab Sample ID: 240-59482-2 Date Sampled: 12/28/2015 1000
Client Matrix: Solid % Moisture: 18.4 Date Received: 12/29/2015 0930

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-212780 | Instrument ID: | A2HP5F |
| Prep Method: | 3540C | Prep Batch: | 240-212549 | Initial Weight/Volume: | 29.97 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 12/31/2015 1744 | | | Injection Volume: | 1 uL |
| Prep Date: | 12/30/2015 0714 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 7.9 | U | 7.9 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 92 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-8**

Lab Sample ID: 240-59482-3 Date Sampled: 12/28/2015 1030
Client Matrix: Solid % Moisture: 16.1 Date Received: 12/29/2015 0930

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8015C | Analysis Batch: | 240-212780 | Instrument ID: | A2HP5F |
| Prep Method: | 3540C | Prep Batch: | 240-212549 | Initial Weight/Volume: | 29.78 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 5 mL |
| Analysis Date: | 12/31/2015 1815 | | | Injection Volume: | 1 uL |
| Prep Date: | 12/30/2015 0714 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (mg/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | | 42 | | 7.7 | 20 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| o-Terphenyl (Surr) | 104 | | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-7**

Lab Sample ID: 240-59482-2 Date Sampled: 12/28/2015 1000
Client Matrix: Solid % Moisture: 18.4 Date Received: 12/29/2015 0930

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-212845 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-212544 | Initial Weight/Volume: | 29.83 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 01/04/2016 1019 | | | Injection Volume: | 1 uL |
| Prep Date: | 12/30/2015 0654 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 15 | U | 15 | 41 |
| Aroclor-1221 | | 20 | U | 20 | 41 |
| Aroclor-1232 | | 25 | U | 25 | 41 |
| Aroclor-1242 | | 14 | U | 14 | 41 |
| Aroclor-1248 | | 9.9 | U | 9.9 | 41 |
| Aroclor-1254 | | 17 | U | 17 | 41 |
| Aroclor-1260 | | 11 | U | 11 | 41 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 84 | | 13 - 134 |
| DCB Decachlorobiphenyl | 72 | | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Client Sample ID:** **F-VB-UST-P-8**

Lab Sample ID: 240-59482-3 Date Sampled: 12/28/2015 1030
Client Matrix: Solid % Moisture: 16.1 Date Received: 12/29/2015 0930

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| | | | | | |
|---|---|---|---|---|---|
| Analysis Method: | 8082A | Analysis Batch: | 240-212845 | Instrument ID: | A2HP12 |
| Prep Method: | 3540C | Prep Batch: | 240-212544 | Initial Weight/Volume: | 29.77 g |
| Dilution: | 1.0 | | | Final Weight/Volume: | 10 mL |
| Analysis Date: | 01/04/2016 1037 | | | Injection Volume: | 1 uL |
| Prep Date: | 12/30/2015 0654 | | | Result Type: | PRIMARY |

| Analyte | DryWt Corrected: Y | Result (ug/Kg) | Qualifier | MDL | RL |
|---|---|---|---|---|---|
| Aroclor-1016 | | 14 | U | 14 | 40 |
| Aroclor-1221 | | 19 | U | 19 | 40 |
| Aroclor-1232 | | 24 | U | 24 | 40 |
| Aroclor-1242 | | 13 | U | 13 | 40 |
| Aroclor-1248 | | 9.6 | U | 9.6 | 40 |
| Aroclor-1254 | | 17 | U | 17 | 40 |
| Aroclor-1260 | | 11 | U | 11 | 40 |

| Surrogate | %Rec | Qualifier | Acceptance Limits |
|---|---|---|---|
| Tetrachloro-m-xylene | 57 | | 13 - 134 |
| DCB Decachlorobiphenyl | 92 | | 10 - 155 |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-P-7**

Lab Sample ID: 240-59482-2 Date Sampled: 12/28/2015 1000
Client Matrix: Solid Date Received: 12/29/2015 0930

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 82 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-212487 | Analysis Date: 12/29/2015 1350 | | | | | | DryWt Corrected: N |
| Percent Moisture | 18 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-212487 | Analysis Date: 12/29/2015 1350 | | | | | | DryWt Corrected: N |

**General Chemistry**

**Client Sample ID:** **F-VB-UST-P-8**

Lab Sample ID: 240-59482-3 Date Sampled: 12/28/2015 1030
Client Matrix: Solid Date Received: 12/29/2015 0930

| Analyte | Result | Qual | Units | MDL | RL | Dil | Method |
|---|---|---|---|---|---|---|---|
| Percent Solids | 84 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-212487 | Analysis Date: 12/29/2015 1350 | | | | | | DryWt Corrected: N |
| Percent Moisture | 16 | | % | 0.10 | 0.10 | 1.0 | Moisture |
| Analysis Batch: 240-212487 | Analysis Date: 12/29/2015 1350 | | | | | | DryWt Corrected: N |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Surrogate Recovery Report

**8260C Volatile Organic Compounds by GC/MS**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | DBFM %Rec | DCA %Rec | TOL %Rec | BFB %Rec |
|---|---|---|---|---|---|
| 240-59482-2 | F-VB-UST-P-7 | 93 | 100 | 91 | 88 |
| 240-59482-3 | F-VB-UST-P-8 | 90 | 97 | 91 | 87 |
| MB 240-212606/24 | | 93 | 97 | 88 | 86 |
| LCS 240-212606/23 | | 94 | 92 | 97 | 91 |
| LCS 240-212606/25 | | 90 | 92 | 93 | 89 |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 95 | 97 | 92 | 91 |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 95 | 100 | 91 | 87 |

| Surrogate | Acceptance Limits |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 37-132 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 58-123 |
| TOL = Toluene-d8 (Surr) | 67-125 |
| BFB = 4-Bromofluorobenzene (Surr) | 52-136 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Surrogate Recovery Report

**8260C Volatile Organic Compounds by GC/MS**

**Client Matrix: Water**

| Lab Sample ID | Client Sample ID | DBFM %Rec | DCA %Rec | TOL %Rec | BFB %Rec |
|---|---|---|---|---|---|
| 240-59482-1 | TB-122815 | 93 | 99 | 102 | 112 |
| MB 240-212893/8 | | 94 | 97 | 102 | 111 |
| LCS 240-212893/5 | | 101 | 96 | 100 | 114 |
| 240-59545-B-2 MS | | 100 | 102 | 103 | 117 |
| 240-59545-B-2 MSD | | 105 | 97 | 98 | 116 |

| Surrogate | Acceptance Limits |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 79-120 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 78-125 |
| TOL = Toluene-d8 (Surr) | 80-120 |
| BFB = 4-Bromofluorobenzene (Surr) | 61-120 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Surrogate Recovery Report

**8270D Semivolatile Organic Compounds (GC/MS)**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | 2FP %Rec | PHL %Rec | NBZ %Rec | FBP %Rec | TBP %Rec | TPH %Rec |
|---|---|---|---|---|---|---|---|
| 240-59482-2 | F-VB-UST-P-7 | 81 | 83 | 80 | 77 | 88 | 87 |
| 240-59482-3 | F-VB-UST-P-8 | 83 | 85 | 81 | 77 | 101 | 87 |
| MB 240-212548/7-A | | 81 | 86 | 79 | 78 | 80 | 89 |
| LCS 240-212548/8-A | | 77 | 80 | 82 | 74 | 90 | 81 |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 78 | 80 | 84 | 77 | 101 | 89 |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 77 | 79 | 86 | 78 | 90 | 89 |

| Surrogate | Acceptance Limits |
|---|---|
| 2FP = 2-Fluorophenol (Surr) | 24-110 |
| PHL = Phenol-d5 (Surr) | 26-110 |
| NBZ = Nitrobenzene-d5 (Surr) | 20-110 |
| FBP = 2-Fluorobiphenyl (Surr) | 24-110 |
| TBP = 2,4,6-Tribromophenol (Surr) | 10-110 |
| TPH = Terphenyl-d14 (Surr) | 36-110 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Surrogate Recovery Report

**8015C Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics)**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | TFT1 %Rec |
|---|---|---|
| 240-59482-2 | F-VB-UST-P-7 | 70 |
| 240-59482-3 | F-VB-UST-P-8 | 67 |
| MB 240-212559/1-A | | 68 |
| LCS 240-212559/2-A | | 90 |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 82 |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 82 |

| Surrogate | Acceptance Limits |
|---|---|
| TFT = Trifluorotoluene (Surr) | 40-139 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Surrogate Recovery Report

**8015C Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | OTPH1 %Rec |
|---|---|---|
| 240-59482-2 | F-VB-UST-P-7 | 92 |
| 240-59482-3 | F-VB-UST-P-8 | 104 |
| MB 240-212549/8-A | | 90 |
| LCS 240-212549/9-A | | 97 |

| Surrogate | Acceptance Limits |
|---|---|
| OTPH = o-Terphenyl (Surr) | 40-160 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Surrogate Recovery Report

**8082A Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

**Client Matrix: Solid**

| Lab Sample ID | Client Sample ID | TCX2 %Rec | DCB2 %Rec |
|---|---|---|---|
| 240-59482-2 | F-VB-UST-P-7 | 84 | 72 |
| 240-59482-3 | F-VB-UST-P-8 | 57 | 92 |
| MB 240-212544/7-A | | 106 | 110 |
| LCS 240-212544/8-A | | 102 | 110 |
| 240-59503-G-4-B MS | | 44 | 43 |
| 240-59503-G-4-C MSD | | 31p | 37 |

| Surrogate | Acceptance Limits |
|---|---|
| TCX = Tetrachloro-m-xylene | 13-134 |
| DCB = DCB Decachlorobiphenyl | 10-155 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Method Blank - Batch: 240-212606** **Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-212606/24 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 147321.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 12/30/2015 2000 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.50 | U | 0.50 | 5.0 |
| 1,1,1-Trichloroethane | 0.65 | U | 0.65 | 5.0 |
| 1,1,2,2-Tetrachloroethane | 0.31 | U | 0.31 | 5.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 0.42 | 5.0 |
| 1,1-Dichloroethane | 0.34 | U | 0.34 | 5.0 |
| 1,1-Dichloroethene | 0.80 | U | 0.80 | 5.0 |
| 1,1-Dichloropropene | 0.37 | U | 0.37 | 5.0 |
| 1,2,3-Trichlorobenzene | 0.562 | J | 0.50 | 5.0 |
| 1,2,3-Trichloropropane | 0.78 | U | 0.78 | 5.0 |
| 1,2,3-Trimethylbenzene | 0.46 | U | 0.46 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.460 | J | 0.38 | 5.0 |
| 1,2,4-Trimethylbenzene | 0.58 | U | 0.58 | 5.0 |
| 1,2-Dibromo-3-Chloropropane | 2.7 | U | 2.7 | 10 |
| 1,2-Dichlorobenzene | 0.34 | U | 0.34 | 5.0 |
| 1,2-Dichloroethane | 0.46 | U | 0.46 | 5.0 |
| 1,2-Dichloropropane | 0.14 | U | 0.14 | 5.0 |
| 1,3-Dichlorobenzene | 0.69 | U | 0.69 | 5.0 |
| 1,3-Dichloropropane | 0.40 | U | 0.40 | 5.0 |
| 1,4-Dichlorobenzene | 0.28 | U | 0.28 | 5.0 |
| 2,2-Dichloropropane | 0.44 | U | 0.44 | 5.0 |
| 2-Butanone (MEK) | 1.1 | U | 1.1 | 20 |
| 2-Chloroethyl vinyl ether | 0.64 | U | 0.64 | 50 |
| 2-Chlorotoluene | 0.30 | U | 0.30 | 5.0 |
| 2-Hexanone | 0.68 | U | 0.68 | 20 |
| 4-Chlorotoluene | 0.26 | U | 0.26 | 5.0 |
| p-Isopropyltoluene | 0.41 | U | 0.41 | 5.0 |
| 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 1.2 | 20 |
| Acetone | 5.7 | U | 5.7 | 20 |
| Benzene | 0.90 | U | 0.90 | 5.0 |
| Bromobenzene | 0.33 | U | 0.33 | 5.0 |
| Bromoform | 0.24 | U | 0.24 | 5.0 |
| Bromomethane | 0.42 | U | 0.42 | 5.0 |
| Carbon disulfide | 0.59 | U | 0.59 | 5.0 |
| Carbon tetrachloride | 0.65 | U | 0.65 | 5.0 |
| Chlorobenzene | 0.53 | U | 0.53 | 5.0 |
| Dibromochloromethane | 0.37 | U | 0.37 | 5.0 |
| Chloroethane | 0.44 | U | 0.44 | 5.0 |
| Chloroform | 0.36 | U | 0.36 | 5.0 |
| Chloromethane | 0.79 | U | 0.79 | 5.0 |
| cis-1,2-Dichloroethene | 0.28 | U | 0.28 | 5.0 |
| cis-1,3-Dichloropropene | 0.71 | U | 0.71 | 5.0 |
| Dibromomethane | 0.42 | U | 0.42 | 5.0 |
| Bromochloromethane | 0.33 | U | 0.33 | 5.0 |
| Dichlorodifluoromethane | 0.28 | U | 0.28 | 5.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 5.0 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Method Blank - Batch: 240-212606**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-212606/24 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 147321.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 12/30/2015 2000 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| tert-Butyl alcohol (TBA) | 5.7 | U | 5.7 | 200 |
| Ethylbenzene | 0.32 | U | 0.32 | 5.0 |
| Hexachlorobutadiene | 0.40 | U | 0.40 | 5.0 |
| Diisopropyl ether | 0.59 | U | 0.59 | 10 |
| Isopropylbenzene | 0.20 | U | 0.20 | 5.0 |
| Methyl tert-butyl ether | 0.43 | U | 0.43 | 5.0 |
| Methylene Chloride | 0.74 | U | 0.74 | 5.0 |
| m-Xylene & p-Xylene | 0.42 | U | 0.42 | 10 |
| Naphthalene | 0.618 | J | 0.35 | 5.0 |
| n-Butylbenzene | 0.59 | U | 0.59 | 5.0 |
| N-Propylbenzene | 0.60 | U | 0.60 | 5.0 |
| o-Xylene | 0.57 | U | 0.57 | 5.0 |
| sec-Butylbenzene | 0.91 | U | 0.91 | 5.0 |
| Styrene | 0.40 | U | 0.40 | 5.0 |
| Tert-amyl methyl ether | 0.43 | U | 0.43 | 5.0 |
| Ethyl t-butyl ether | 0.41 | U | 0.41 | 5.0 |
| Tetrachloroethene | 0.80 | U | 0.80 | 5.0 |
| Toluene | 0.27 | U | 0.27 | 5.0 |
| trans-1,2-Dichloroethene | 0.32 | U | 0.32 | 5.0 |
| trans-1,3-Dichloropropene | 0.29 | U | 0.29 | 5.0 |
| Trichloroethene | 0.38 | U | 0.38 | 5.0 |
| Trichlorofluoromethane | 0.29 | U | 0.29 | 5.0 |
| Vinyl acetate | 0.66 | U | 0.66 | 10 |
| Vinyl chloride | 0.30 | U | 0.30 | 5.0 |
| Xylenes, Total | 0.54 | U | 0.54 | 10 |
| tert-Butylbenzene | 0.67 | U | 0.67 | 5.0 |
| 1,2-Dibromoethane | 0.53 | U | 0.53 | 5.0 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 88 | 67 - 125 |
| Dibromofluoromethane (Surr) | 93 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 86 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 97 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Lab Control Sample - Batch: 240-212606**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-212606/23 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 147320.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 12/30/2015 1939 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 50.0 | 58.6 | 117 | 80 - 120 | |
| 1,1,1-Trichloroethane | 50.0 | 54.0 | 108 | 77 - 126 | |
| 1,1,2,2-Tetrachloroethane | 50.0 | 53.5 | 107 | 77 - 123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 50.0 | 45.3 | 91 | 80 - 138 | |
| 1,1-Dichloroethane | 50.0 | 51.1 | 102 | 76 - 120 | |
| 1,1-Dichloroethene | 50.0 | 42.2 | 84 | 75 - 135 | |
| 1,1-Dichloropropene | 50.0 | 54.6 | 109 | 80 - 120 | |
| 1,2,3-Trichlorobenzene | 50.0 | 45.6 | 91 | 61 - 121 | |
| 1,2,3-Trichloropropane | 50.0 | 57.2 | 114 | 80 - 126 | |
| 1,2,4-Trichlorobenzene | 50.0 | 45.6 | 91 | 64 - 124 | |
| 1,2,4-Trimethylbenzene | 50.0 | 52.6 | 105 | 80 - 129 | |
| 1,2-Dibromo-3-Chloropropane | 50.0 | 55.6 | 111 | 61 - 132 | |
| 1,2-Dichlorobenzene | 50.0 | 48.1 | 96 | 76 - 120 | |
| 1,2-Dichloroethane | 50.0 | 52.5 | 105 | 72 - 120 | |
| 1,2-Dichloropropane | 50.0 | 52.7 | 105 | 80 - 120 | |
| 1,3-Dichlorobenzene | 50.0 | 47.0 | 94 | 78 - 120 | |
| 1,3-Dichloropropane | 50.0 | 51.7 | 103 | 80 - 120 | |
| 1,4-Dichlorobenzene | 50.0 | 47.3 | 95 | 75 - 120 | |
| 2,2-Dichloropropane | 50.0 | 54.3 | 109 | 69 - 135 | |
| 2-Butanone (MEK) | 100 | 96.3 | 96 | 52 - 131 | |
| 2-Chloroethyl vinyl ether | 50.0 | 54.3 | 109 | 28 - 173 | |
| 2-Chlorotoluene | 50.0 | 52.2 | 104 | 78 - 120 | |
| 2-Hexanone | 100 | 99.2 | 99 | 64 - 136 | |
| 4-Chlorotoluene | 50.0 | 48.3 | 97 | 79 - 120 | |
| p-Isopropyltoluene | 50.0 | 53.4 | 107 | 77 - 131 | |
| 4-Methyl-2-pentanone (MIBK) | 100 | 109 | 109 | 67 - 135 | |
| Acetone | 100 | 94.9 | 95 | 41 - 137 | |
| Benzene | 50.0 | 50.8 | 102 | 79 - 120 | |
| Bromobenzene | 50.0 | 51.6 | 103 | 80 - 120 | |
| Bromoform | 50.0 | 54.9 | 110 | 62 - 133 | |
| Bromomethane | 20.0 | 19.2 | 96 | 42 - 136 | |
| Carbon disulfide | 50.0 | 54.9 | 110 | 62 - 146 | |
| Carbon tetrachloride | 50.0 | 57.7 | 115 | 71 - 129 | |
| Chlorobenzene | 50.0 | 50.5 | 101 | 78 - 120 | |
| Dibromochloromethane | 50.0 | 54.7 | 109 | 72 - 127 | |
| Chloroethane | 20.0 | 19.5 | 98 | 58 - 120 | |
| Chloroform | 50.0 | 51.8 | 104 | 77 - 120 | |
| Chloromethane | 20.0 | 20.5 | 102 | 50 - 120 | |
| cis-1,2-Dichloroethene | 50.0 | 52.6 | 105 | 76 - 120 | |
| cis-1,3-Dichloropropene | 50.0 | 53.5 | 107 | 74 - 128 | |
| Dibromomethane | 50.0 | 53.8 | 108 | 80 - 120 | |
| Bromochloromethane | 50.0 | 52.0 | 104 | 79 - 120 | |
| Dichlorodifluoromethane | 20.0 | 20.7 | 103 | 26 - 120 | |
| Bromodichloromethane | 50.0 | 54.4 | 109 | 80 - 122 | |
| tert-Butyl alcohol (TBA) | 500 | 450 | 90 | 60 - 125 | |
| Ethylbenzene | 50.0 | 51.6 | 103 | 79 - 120 | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Lab Control Sample - Batch: 240-212606**

**Method: 8260C**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-212606/23 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | 147320.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 12/30/2015 1939 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Hexachlorobutadiene | 50.0 | 51.3 | 103 | 54 - 131 | |
| Isopropylbenzene | 50.0 | 54.4 | 109 | 76 - 122 | |
| Methyl tert-butyl ether | 50.0 | 53.0 | 106 | 49 - 165 | |
| Methylene Chloride | 50.0 | 50.7 | 101 | 75 - 120 | |
| m-Xylene & p-Xylene | 50.0 | 53.8 | 108 | 80 - 120 | |
| Naphthalene | 50.0 | 51.1 | 102 | 65 - 123 | |
| n-Butylbenzene | 50.0 | 50.3 | 101 | 68 - 135 | |
| N-Propylbenzene | 50.0 | 53.2 | 106 | 80 - 129 | |
| o-Xylene | 50.0 | 54.6 | 109 | 80 - 120 | |
| sec-Butylbenzene | 50.0 | 55.3 | 111 | 74 - 129 | |
| Styrene | 50.0 | 53.8 | 108 | 80 - 120 | |
| Tetrachloroethene | 50.0 | 51.4 | 103 | 79 - 120 | |
| Toluene | 50.0 | 53.2 | 106 | 75 - 120 | |
| trans-1,2-Dichloroethene | 50.0 | 53.2 | 106 | 78 - 120 | |
| trans-1,3-Dichloropropene | 50.0 | 55.5 | 111 | 73 - 131 | |
| Trichloroethene | 50.0 | 53.1 | 106 | 79 - 120 | |
| Trichlorofluoromethane | 20.0 | 21.1 | 106 | 57 - 146 | |
| Vinyl acetate | 50.0 | 48.6 | 97 | 70 - 130 | |
| Vinyl chloride | 20.0 | 20.2 | 101 | 57 - 120 | |
| Xylenes, Total | 100 | 108 | 108 | 80 - 120 | |
| tert-Butylbenzene | 50.0 | 54.9 | 110 | 76 - 126 | |
| 1,2-Dibromoethane | 50.0 | 56.1 | 112 | 80 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 97 | 67 - 125 |
| Dibromofluoromethane (Surr) | 94 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 91 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 92 | 58 - 123 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 93 | 67 - 125 |
| Dibromofluoromethane (Surr) | 90 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 89 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 92 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212657**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-212657 | Lab File ID: | 147327.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.65 g |
| Analysis Date: | 12/30/2015 2209 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-212657 | Lab File ID: | 147328.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.47 g |
| Analysis Date: | 12/30/2015 2231 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1,1,2-Tetrachloroethane | 82 | 74 | 34 - 135 | 7 | 30 | | |
| 1,1,1-Trichloroethane | 93 | 98 | 51 - 128 | 8 | 30 | | |
| 1,1,2,2-Tetrachloroethane | 80 | 85 | 16 - 179 | 9 | 30 | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 96 | 104 | 50 - 147 | 11 | 30 | | |
| 1,1-Dichloroethane | 85 | 91 | 54 - 122 | 10 | 30 | | |
| 1,1-Dichloroethene | 89 | 102 | 49 - 157 | 17 | 30 | | |
| 1,1-Dichloropropene | 93 | 93 | 50 - 122 | 3 | 30 | | |
| 1,2,3-Trichlorobenzene | 42 | 34 | 10 - 120 | 18 | 30 | | |
| 1,2,3-Trichloropropane | 89 | 90 | 32 - 174 | 5 | 30 | | |
| 1,2,4-Trichlorobenzene | 43 | 31 | 10 - 120 | 29 | 30 | | |
| 1,2,4-Trimethylbenzene | 65 | 45 | 10 - 173 | 30 | 30 | | |
| 1,2-Dibromo-3-Chloropropane | 74 | 87 | 10 - 153 | 20 | 30 | | |
| 1,2-Dichlorobenzene | 57 | 42 | 17 - 122 | 27 | 30 | | |
| 1,2-Dichloroethane | 82 | 89 | 49 - 123 | 11 | 30 | | |
| 1,2-Dichloropropane | 87 | 87 | 61 - 120 | 3 | 30 | | |
| 1,3-Dichlorobenzene | 55 | 37 | 16 - 126 | 36 | 30 | | F2 |
| 1,3-Dichloropropane | 76 | 78 | 54 - 128 | 5 | 30 | | |
| 1,4-Dichlorobenzene | 51 | 36 | 15 - 121 | 32 | 30 | | F2 |
| 2,2-Dichloropropane | 85 | 90 | 49 - 132 | 9 | 30 | | |
| 2-Butanone (MEK) | 78 | 91 | 30 - 143 | 18 | 30 | | |
| 2-Chloroethyl vinyl ether | 61 | 71 | 10 - 162 | 19 | 30 | J | J |
| 2-Chlorotoluene | 65 | 43 | 11 - 162 | 38 | 30 | | F2 |
| 2-Hexanone | 83 | 92 | 37 - 147 | 13 | 30 | | |
| 4-Chlorotoluene | 59 | 38 | 15 - 149 | 40 | 30 | | F2 |
| p-Isopropyltoluene | 63 | 36 | 10 - 165 | 52 | 30 | | F2 |
| 4-Methyl-2-pentanone (MIBK) | 90 | 104 | 43 - 147 | 19 | 30 | | |
| Acetone | 61 | 80 | 24 - 140 | 20 | 30 | | |
| Benzene | 85 | 85 | 53 - 120 | 4 | 30 | | |
| Bromobenzene | 65 | 51 | 24 - 144 | 22 | 30 | | |
| Bromoform | 78 | 76 | 18 - 129 | 1 | 30 | | |
| Bromomethane | 74 | 80 | 33 - 130 | 11 | 30 | | |
| Carbon disulfide | 80 | 82 | 20 - 151 | 5 | 30 | | |
| Carbon tetrachloride | 100 | 101 | 32 - 137 | 5 | 30 | | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212657**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-212657 | Lab File ID: | 147327.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.65 g |
| Analysis Date: | 12/30/2015 2209 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-212657 | Lab File ID: | 147328.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.47 g |
| Analysis Date: | 12/30/2015 2231 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Chlorobenzene | 71 | 56 | 37 - 120 | 19 | 30 | | |
| Dibromochloromethane | 78 | 78 | 29 - 135 | 3 | 30 | | |
| Chloroethane | 79 | 85 | 45 - 120 | 10 | 30 | | |
| Chloroform | 83 | 87 | 53 - 120 | 8 | 30 | | |
| Chloromethane | 85 | 101 | 34 - 120 | 20 | 30 | | |
| cis-1,2-Dichloroethene | 80 | 85 | 50 - 120 | 10 | 30 | | |
| cis-1,3-Dichloropropene | 78 | 75 | 27 - 133 | 2 | 30 | | |
| Dibromomethane | 80 | 83 | 52 - 123 | 7 | 30 | | |
| Bromochloromethane | 77 | 84 | 53 - 120 | 12 | 30 | | |
| Dichlorodifluoromethane | 95 | 108 | 17 - 120 | 16 | 30 | | |
| Bromodichloromethane | 85 | 84 | 35 - 132 | 2 | 30 | | |
| tert-Butyl alcohol (TBA) | 84 | 97 | 60 - 125 | 17 | 30 | | |
| Ethylbenzene | 73 | 55 | 30 - 131 | 25 | 30 | | |
| Hexachlorobutadiene | 45 | 22 | 10 - 131 | 66 | 30 | | F2 |
| Isopropylbenzene | 72 | 51 | 21 - 134 | 32 | 30 | | F2 |
| Methyl tert-butyl ether | 85 | 95 | 51 - 157 | 13 | 30 | | |
| Methylene Chloride | 75 | 81 | 54 - 120 | 11 | 30 | | |
| m-Xylene & p-Xylene | 72 | 57 | 29 - 131 | 19 | 30 | | |
| Naphthalene | 56 | 55 | 10 - 124 | 1 | 30 | | |
| n-Butylbenzene | 54 | 28 | 10 - 148 | 62 | 30 | | F2 |
| N-Propylbenzene | 68 | 42 | 10 - 178 | 45 | 30 | | F2 |
| o-Xylene | 76 | 58 | 29 - 134 | 22 | 30 | | |
| sec-Butylbenzene | 66 | 38 | 10 - 172 | 50 | 30 | | F2 |
| Styrene | 69 | 50 | 27 - 127 | 28 | 30 | | |
| Tetrachloroethene | 78 | 62 | 31 - 135 | 20 | 30 | | |
| Toluene | 80 | 69 | 39 - 129 | 11 | 30 | | |
| trans-1,2-Dichloroethene | 81 | 86 | 50 - 123 | 9 | 30 | | |
| trans-1,3-Dichloropropene | 69 | 69 | 28 - 137 | 4 | 30 | | |
| Trichloroethene | 82 | 79 | 10 - 177 | 1 | 30 | | |
| Trichlorofluoromethane | 96 | 110 | 36 - 142 | 17 | 30 | | |
| Vinyl acetate | 61 | 78 | 70 - 130 | 28 | 30 | F1 | |
| Vinyl chloride | 84 | 97 | 42 - 120 | 18 | 30 | | |
| Xylenes, Total | 74 | 58 | 30 - 131 | 20 | 30 | | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212657**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-212657 | Lab File ID: | 147327.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.65 g |
| Analysis Date: | 12/30/2015 2209 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212606 | Instrument ID: | A3UX14 |
| Client Matrix: | Solid | Prep Batch: | 240-212657 | Lab File ID: | 147328.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.47 g |
| Analysis Date: | 12/30/2015 2231 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 1425 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| tert-Butylbenzene | 70 | 45 | 10 - 163 | 40 | 30 | | F2 |
| 1,2-Dibromoethane | 78 | 78 | 45 - 127 | 3 | 30 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 92 | 91 | 67 - 125 |
| Dibromofluoromethane (Surr) | 95 | 95 | 37 - 132 |
| 4-Bromofluorobenzene (Surr) | 91 | 87 | 52 - 136 |
| 1,2-Dichloroethane-d4 (Surr) | 97 | 100 | 58 - 123 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212657**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Units: ug/Kg | MSD Lab Sample ID: | 240-59482-2 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 12/30/2015 2209 | | Analysis Date: | 12/30/2015 2231 | |
| Prep Date: | 12/30/2015 1425 | | Prep Date: | 12/30/2015 1425 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.58 | U | 54.2 | 56.0 | 44.5 | | 41.7 | |
| 1,1,1-Trichloroethane | 0.76 | U | 54.2 | 56.0 | 50.6 | | 55.0 | |
| 1,1,2,2-Tetrachloroethane | 0.36 | U | 54.2 | 56.0 | 43.6 | | 47.6 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 54.2 | 56.0 | 52.2 | | 58.3 | |
| 1,1-Dichloroethane | 0.40 | U | 54.2 | 56.0 | 45.9 | | 50.9 | |
| 1,1-Dichloroethene | 0.93 | U | 54.2 | 56.0 | 48.2 | | 56.9 | |
| 1,1-Dichloropropene | 0.43 | U | 54.2 | 56.0 | 50.4 | | 51.9 | |
| 1,2,3-Trichlorobenzene | 0.58 | U | 54.2 | 56.0 | 22.8 | | 19.0 | |
| 1,2,3-Trichloropropane | 0.91 | U | 54.2 | 56.0 | 48.1 | | 50.5 | |
| 1,2,4-Trichlorobenzene | 0.44 | U | 54.2 | 56.0 | 23.3 | | 17.4 | |
| 1,2,4-Trimethylbenzene | 1.9 | J | 54.2 | 56.0 | 37.0 | | 27.3 | |
| 1,2-Dibromo-3-Chloropropane | 3.2 | U | 54.2 | 56.0 | 40.1 | | 48.9 | |
| 1,2-Dichlorobenzene | 0.40 | U | 54.2 | 56.0 | 30.9 | | 23.5 | |
| 1,2-Dichloroethane | 0.54 | U | 54.2 | 56.0 | 44.3 | | 49.6 | |
| 1,2-Dichloropropane | 0.16 | U | 54.2 | 56.0 | 47.0 | | 48.5 | |
| 1,3-Dichlorobenzene | 0.80 | U | 54.2 | 56.0 | 29.9 | | 20.8 | F2 |
| 1,3-Dichloropropane | 0.47 | U | 54.2 | 56.0 | 41.4 | | 43.5 | |
| 1,4-Dichlorobenzene | 0.33 | U | 54.2 | 56.0 | 27.7 | | 20.0 | F2 |
| 2,2-Dichloropropane | 0.51 | U | 54.2 | 56.0 | 46.0 | | 50.4 | |
| 2-Butanone (MEK) | 4.8 | J | 108 | 112 | 89.2 | | 106 | |
| 2-Chloroethyl vinyl ether | 0.75 | U | 54.2 | 56.0 | 33.0 | J | 39.9 | J |
| 2-Chlorotoluene | 0.35 | U | 54.2 | 56.0 | 35.4 | | 24.1 | F2 |
| 2-Hexanone | 0.79 | U | 108 | 112 | 90.1 | | 103 | |
| 4-Chlorotoluene | 0.30 | U | 54.2 | 56.0 | 32.1 | | 21.3 | F2 |
| p-Isopropyltoluene | 0.48 | U | 54.2 | 56.0 | 34.2 | | 20.0 | F2 |
| 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 108 | 112 | 97.1 | | 117 | |
| Acetone | 36 | | 108 | 112 | 102 | | 125 | |
| Benzene | 1.0 | U | 54.2 | 56.0 | 45.9 | | 47.6 | |
| Bromobenzene | 0.38 | U | 54.2 | 56.0 | 35.4 | | 28.4 | |
| Bromoform | 0.28 | U | 54.2 | 56.0 | 42.3 | | 42.8 | |
| Bromomethane | 0.49 | U | 21.7 | 22.4 | 16.0 | | 17.9 | |
| Carbon disulfide | 0.69 | U | 54.2 | 56.0 | 43.4 | | 45.7 | |
| Carbon tetrachloride | 0.76 | U | 54.2 | 56.0 | 54.2 | | 56.7 | |
| Chlorobenzene | 0.62 | U | 54.2 | 56.0 | 38.3 | | 31.6 | |
| Dibromochloromethane | 0.43 | U | 54.2 | 56.0 | 42.4 | | 43.6 | |
| Chloroethane | 0.51 | U | 21.7 | 22.4 | 17.2 | | 19.1 | |
| Chloroform | 0.42 | U | 54.2 | 56.0 | 45.3 | | 49.0 | |
| Chloromethane | 0.92 | U | 21.7 | 22.4 | 18.5 | | 22.6 | |
| cis-1,2-Dichloroethene | 0.33 | U | 54.2 | 56.0 | 43.2 | | 47.7 | |
| cis-1,3-Dichloropropene | 0.83 | U | 54.2 | 56.0 | 42.5 | | 41.8 | |
| Dibromomethane | 0.49 | U | 54.2 | 56.0 | 43.3 | | 46.2 | |
| Bromochloromethane | 0.38 | U | 54.2 | 56.0 | 41.7 | | 46.8 | |
| Dichlorodifluoromethane | 0.33 | U | 21.7 | 22.4 | 20.5 | | 24.1 | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212657**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Units: ug/Kg | MSD Lab Sample ID: | 240-59482-2 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 12/30/2015 2209 | | Analysis Date: | 12/30/2015 2231 | |
| Prep Date: | 12/30/2015 1425 | | Prep Date: | 12/30/2015 1425 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| Bromodichloromethane | 0.34 | U | 54.2 | 56.0 | 46.2 | | 47.1 | |
| tert-Butyl alcohol (TBA) | 6.7 | U | 542 | 560 | 454 | | 541 | |
| Ethylbenzene | 0.66 | J | 54.2 | 56.0 | 40.5 | | 31.4 | |
| Hexachlorobutadiene | 0.47 | U | 54.2 | 56.0 | 24.2 | | 12.2 | F2 |
| Isopropylbenzene | 0.24 | J | 54.2 | 56.0 | 39.5 | | 28.6 | F2 |
| Methyl tert-butyl ether | 0.50 | U | 54.2 | 56.0 | 46.4 | | 53.0 | |
| Methylene Chloride | 0.86 | U | 54.2 | 56.0 | 40.6 | | 45.3 | |
| m-Xylene & p-Xylene | 2.9 | J | 54.2 | 56.0 | 42.1 | | 34.9 | |
| Naphthalene | 0.61 | J | 54.2 | 56.0 | 31.1 | | 31.3 | |
| n-Butylbenzene | 0.69 | U | 54.2 | 56.0 | 29.5 | | 15.6 | F2 |
| N-Propylbenzene | 0.70 | U | 54.2 | 56.0 | 36.8 | | 23.4 | F2 |
| o-Xylene | 1.2 | J | 54.2 | 56.0 | 42.2 | | 33.9 | |
| sec-Butylbenzene | 1.1 | U | 54.2 | 56.0 | 35.7 | | 21.4 | F2 |
| Styrene | 0.47 | U | 54.2 | 56.0 | 37.2 | | 28.1 | |
| Tetrachloroethene | 0.93 | U | 54.2 | 56.0 | 42.6 | | 34.7 | |
| Toluene | 1.6 | J | 54.2 | 56.0 | 45.1 | | 40.2 | |
| trans-1,2-Dichloroethene | 0.37 | U | 54.2 | 56.0 | 44.0 | | 48.3 | |
| trans-1,3-Dichloropropene | 0.34 | U | 54.2 | 56.0 | 37.3 | | 38.8 | |
| Trichloroethene | 0.44 | U | 54.2 | 56.0 | 44.7 | | 44.3 | |
| Trichlorofluoromethane | 0.34 | U | 21.7 | 22.4 | 20.8 | | 24.7 | |
| Vinyl acetate | 0.77 | U | 54.2 | 56.0 | 32.9 | F1 | 43.6 | |
| Vinyl chloride | 0.35 | U | 21.7 | 22.4 | 18.3 | | 21.8 | |
| Xylenes, Total | 4.1 | J | 108 | 112 | 84.3 | | 68.8 | |
| tert-Butylbenzene | 0.78 | U | 54.2 | 56.0 | 38.2 | | 25.3 | F2 |
| 1,2-Dibromoethane | 0.62 | U | 54.2 | 56.0 | 42.6 | | 43.9 | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Method Blank - Batch: 240-212893**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-212893/8 | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC5596.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1126 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1126 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 0.28 | U | 0.28 | 1.0 |
| 1,1,1-Trichloroethane | 0.44 | U | 0.44 | 1.0 |
| 1,1,2,2-Tetrachloroethane | 0.22 | U | 0.22 | 1.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloroethane | 0.30 | U | 0.30 | 1.0 |
| 1,1-Dichloroethene | 0.45 | U | 0.45 | 1.0 |
| 1,1-Dichloropropene | 0.42 | U | 0.42 | 1.0 |
| 1,2,3-Trichlorobenzene | 0.37 | U | 0.37 | 1.0 |
| 1,2,3-Trichloropropane | 0.44 | U | 0.44 | 1.0 |
| 1,2,3-Trimethylbenzene | 0.47 | U | 0.47 | 5.0 |
| 1,2,4-Trichlorobenzene | 0.32 | U | 0.32 | 1.0 |
| 1,2,4-Trimethylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromo-3-Chloropropane | 0.82 | U | 0.82 | 2.0 |
| 1,2-Dichlorobenzene | 0.25 | U | 0.25 | 1.0 |
| 1,2-Dichloroethane | 0.23 | U | 0.23 | 1.0 |
| 1,2-Dichloropropane | 0.25 | U | 0.25 | 1.0 |
| 1,3-Dichlorobenzene | 0.19 | U | 0.19 | 1.0 |
| 1,3-Dichloropropane | 0.19 | U | 0.19 | 1.0 |
| 1,4-Dichlorobenzene | 0.27 | U | 0.27 | 1.0 |
| 2,2-Dichloropropane | 0.26 | U | 0.26 | 1.0 |
| 2-Butanone (MEK) | 1.34 | J | 0.53 | 10 |
| 2-Chloroethyl vinyl ether | 0.59 | U | 0.59 | 10 |
| 2-Chlorotoluene | 0.40 | U | 0.40 | 1.0 |
| 2-Hexanone | 0.48 | U | 0.48 | 10 |
| 4-Chlorotoluene | 0.29 | U | 0.29 | 1.0 |
| p-Isopropyltoluene | 0.43 | U | 0.43 | 1.0 |
| 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 0.99 | 10 |
| Acetone | 0.94 | U | 0.94 | 10 |
| Benzene | 0.35 | U | 0.35 | 1.0 |
| Bromobenzene | 0.35 | U | 0.35 | 1.0 |
| Bromoform | 0.56 | U | 0.56 | 1.0 |
| Bromomethane | 0.44 | U | 0.44 | 1.0 |
| Carbon disulfide | 0.38 | U | 0.38 | 1.0 |
| Carbon tetrachloride | 0.43 | U | 0.43 | 1.0 |
| Chlorobenzene | 0.25 | U | 0.25 | 1.0 |
| Dibromochloromethane | 0.43 | U | 0.43 | 1.0 |
| Chloroethane | 0.32 | U | 0.32 | 1.0 |
| Chloroform | 0.25 | U | 0.25 | 1.0 |
| Chloromethane | 0.44 | U | 0.44 | 1.0 |
| cis-1,2-Dichloroethene | 0.26 | U | 0.26 | 1.0 |
| cis-1,3-Dichloropropene | 0.46 | U | 0.46 | 1.0 |
| Dibromomethane | 0.42 | U | 0.42 | 1.0 |
| Bromochloromethane | 0.50 | U | 0.50 | 1.0 |
| Dichlorodifluoromethane | 0.32 | U | 0.32 | 1.0 |
| Bromodichloromethane | 0.29 | U | 0.29 | 1.0 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Method Blank - Batch: 240-212893**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-212893/8 | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC5596.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1126 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1126 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| tert-Butyl alcohol (TBA) | 4.9 | U | 4.9 | 50 |
| Ethylbenzene | 0.25 | U | 0.25 | 1.0 |
| Hexachlorobutadiene | 0.35 | U | 0.35 | 1.0 |
| Diisopropyl ether | 0.50 | U | 0.50 | 10 |
| Isopropylbenzene | 0.35 | U | 0.35 | 1.0 |
| Methyl tert-butyl ether | 0.20 | U | 0.20 | 1.0 |
| Methylene Chloride | 0.33 | U | 0.33 | 1.0 |
| m-Xylene & p-Xylene | 0.24 | U | 0.24 | 2.0 |
| Naphthalene | 0.45 | U | 0.45 | 1.0 |
| n-Butylbenzene | 0.31 | U | 0.31 | 1.0 |
| N-Propylbenzene | 0.40 | U | 0.40 | 1.0 |
| o-Xylene | 0.25 | U | 0.25 | 1.0 |
| sec-Butylbenzene | 0.48 | U | 0.48 | 1.0 |
| Styrene | 0.45 | U | 0.45 | 1.0 |
| Tert-amyl methyl ether | 0.30 | U | 0.30 | 5.0 |
| Ethyl t-butyl ether | 0.23 | U | 0.23 | 5.0 |
| Tetrachloroethene | 0.31 | U | 0.31 | 1.0 |
| Toluene | 0.23 | U | 0.23 | 1.0 |
| trans-1,2-Dichloroethene | 0.30 | U | 0.30 | 1.0 |
| trans-1,3-Dichloropropene | 0.56 | U | 0.56 | 1.0 |
| Trichloroethene | 0.22 | U | 0.22 | 1.0 |
| Trichlorofluoromethane | 0.49 | U | 0.49 | 1.0 |
| Vinyl acetate | 0.41 | U | 0.41 | 2.0 |
| Vinyl chloride | 0.29 | U | 0.29 | 1.0 |
| Xylenes, Total | 0.52 | U | 0.52 | 2.0 |
| tert-Butylbenzene | 0.41 | U | 0.41 | 1.0 |
| 1,2-Dibromoethane | 0.32 | U | 0.32 | 1.0 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 102 | 80 - 120 |
| Dibromofluoromethane (Surr) | 94 | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 111 | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 97 | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Lab Control Sample - Batch: 240-212893**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-212893/5 | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC5593.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1018 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1018 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 10.0 | 8.23 | 82 | 80 - 120 | |
| 1,1,1-Trichloroethane | 10.0 | 8.79 | 88 | 77 - 123 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 9.40 | 94 | 71 - 123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 10.0 | 10.5 | 105 | 67 - 138 | |
| 1,1-Dichloroethane | 10.0 | 9.36 | 94 | 79 - 125 | |
| 1,1-Dichloroethene | 10.0 | 9.64 | 96 | 76 - 124 | |
| 1,1-Dichloropropene | 10.0 | 10.1 | 101 | 80 - 128 | |
| 1,2,3-Trichlorobenzene | 10.0 | 8.36 | 84 | 66 - 120 | |
| 1,2,3-Trichloropropane | 10.0 | 9.73 | 97 | 72 - 126 | |
| 1,2,4-Trichlorobenzene | 10.0 | 8.53 | 85 | 61 - 120 | |
| 1,2,4-Trimethylbenzene | 10.0 | 9.27 | 93 | 76 - 120 | |
| 1,2-Dibromo-3-Chloropropane | 10.0 | 8.17 | 82 | 50 - 132 | |
| 1,2-Dichlorobenzene | 10.0 | 9.02 | 90 | 79 - 120 | |
| 1,2-Dichloroethane | 10.0 | 9.32 | 93 | 80 - 120 | |
| 1,2-Dichloropropane | 10.0 | 9.80 | 98 | 78 - 124 | |
| 1,3-Dichlorobenzene | 10.0 | 9.06 | 91 | 79 - 120 | |
| 1,3-Dichloropropane | 10.0 | 9.36 | 94 | 80 - 120 | |
| 1,4-Dichlorobenzene | 10.0 | 9.07 | 91 | 79 - 120 | |
| 2,2-Dichloropropane | 10.0 | 7.92 | 79 | 71 - 122 | |
| 2-Butanone (MEK) | 20.0 | 21.5 | 108 | 56 - 138 | |
| 2-Chloroethyl vinyl ether | 10.0 | 9.80 | 98 | 54 - 139 | J |
| 2-Chlorotoluene | 10.0 | 9.14 | 91 | 80 - 120 | |
| 2-Hexanone | 20.0 | 18.5 | 92 | 55 - 141 | |
| 4-Chlorotoluene | 10.0 | 9.06 | 91 | 80 - 120 | |
| p-Isopropyltoluene | 10.0 | 9.23 | 92 | 77 - 120 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 19.6 | 98 | 64 - 135 | |
| Acetone | 20.0 | 19.1 | 95 | 34 - 148 | |
| Benzene | 10.0 | 9.39 | 94 | 80 - 120 | |
| Bromobenzene | 10.0 | 9.14 | 91 | 80 - 120 | |
| Bromoform | 10.0 | 6.29 | 63 | 56 - 122 | |
| Bromomethane | 10.0 | 6.43 | 64 | 38 - 132 | |
| Carbon disulfide | 10.0 | 9.43 | 94 | 65 - 144 | |
| Carbon tetrachloride | 10.0 | 7.89 | 79 | 77 - 131 | |
| Chlorobenzene | 10.0 | 9.32 | 93 | 80 - 120 | |
| Dibromochloromethane | 10.0 | 7.97 | 80 | 74 - 120 | |
| Chloroethane | 10.0 | 7.54 | 75 | 36 - 126 | |
| Chloroform | 10.0 | 9.57 | 96 | 80 - 120 | |
| Chloromethane | 10.0 | 9.55 | 96 | 48 - 133 | |
| cis-1,2-Dichloroethene | 10.0 | 9.75 | 98 | 79 - 120 | |
| cis-1,3-Dichloropropene | 10.0 | 9.84 | 98 | 74 - 126 | |
| Dibromomethane | 10.0 | 9.81 | 98 | 80 - 120 | |
| Bromochloromethane | 10.0 | 9.83 | 98 | 80 - 120 | |
| Dichlorodifluoromethane | 10.0 | 7.57 | 76 | 23 - 136 | |
| Bromodichloromethane | 10.0 | 8.91 | 89 | 80 - 120 | |
| tert-Butyl alcohol (TBA) | 100 | 79.5 | 79 | 38 - 144 | |
| Ethylbenzene | 10.0 | 9.45 | 95 | 80 - 120 | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Lab Control Sample - Batch: 240-212893**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-212893/5 | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC5593.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1018 | Units: | ug/L | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1018 | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Hexachlorobutadiene | 10.0 | 7.63 | 76 | 62 - 120 | |
| Isopropylbenzene | 10.0 | 9.47 | 95 | 77 - 120 | |
| Methyl tert-butyl ether | 10.0 | 9.50 | 95 | 69 - 121 | |
| Methylene Chloride | 10.0 | 8.56 | 86 | 77 - 129 | |
| m-Xylene & p-Xylene | 10.0 | 9.38 | 94 | 80 - 120 | |
| Naphthalene | 10.0 | 9.11 | 91 | 53 - 121 | |
| n-Butylbenzene | 10.0 | 8.99 | 90 | 74 - 120 | |
| N-Propylbenzene | 10.0 | 9.29 | 93 | 79 - 123 | |
| o-Xylene | 10.0 | 9.59 | 96 | 80 - 120 | |
| sec-Butylbenzene | 10.0 | 9.27 | 93 | 77 - 120 | |
| Styrene | 10.0 | 9.62 | 96 | 76 - 122 | |
| Tetrachloroethene | 10.0 | 9.39 | 94 | 78 - 121 | |
| Toluene | 10.0 | 9.29 | 93 | 80 - 120 | |
| trans-1,2-Dichloroethene | 10.0 | 9.74 | 97 | 80 - 124 | |
| trans-1,3-Dichloropropene | 10.0 | 9.12 | 91 | 75 - 131 | |
| Trichloroethene | 10.0 | 9.82 | 98 | 80 - 121 | |
| Trichlorofluoromethane | 10.0 | 9.18 | 92 | 61 - 133 | |
| Vinyl acetate | 10.0 | 8.16 | 82 | 27 - 160 | |
| Vinyl chloride | 10.0 | 9.13 | 91 | 52 - 121 | |
| Xylenes, Total | 20.0 | 19.0 | 95 | 80 - 120 | |
| tert-Butylbenzene | 10.0 | 9.36 | 94 | 77 - 121 | |
| 1,2-Dibromoethane | 10.0 | 9.65 | 96 | 80 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Toluene-d8 (Surr) | 100 | 80 - 120 |
| Dibromofluoromethane (Surr) | 101 | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 114 | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 96 | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212893**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59545-B-2 MS | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC5610.D |
| Dilution: | 500 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1700 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1700 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-59545-B-2 MSD | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC5611.D |
| Dilution: | 500 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1723 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1723 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1,1-Trichloroethane | 85 | 90 | 69 - 122 | 5 | 14 | | |
| 1,1,2,2-Tetrachloroethane | 99 | 99 | 61 - 130 | 0 | 18 | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 88 | 104 | 44 - 140 | 16 | 35 | | |
| 1,1-Dichloroethane | 100 | 106 | 73 - 124 | 3 | 14 | | |
| 1,1-Dichloroethene | 95 | 100 | 67 - 124 | 6 | 24 | | |
| 1,2,4-Trichlorobenzene | 80 | 86 | 48 - 120 | 7 | 28 | | |
| 1,2-Dibromo-3-Chloropropane | 78 | 77 | 42 - 130 | 1 | 24 | | |
| 1,2-Dichlorobenzene | 88 | 92 | 67 - 118 | 4 | 15 | | |
| 1,2-Dichloroethane | 99 | 99 | 74 - 125 | 0 | 24 | | |
| 1,2-Dichloropropane | 97 | 98 | 73 - 122 | 1 | 15 | | |
| 1,3-Dichlorobenzene | 89 | 91 | 65 - 120 | 2 | 15 | | |
| 1,4-Dichlorobenzene | 89 | 92 | 66 - 120 | 2 | 16 | | |
| 2-Butanone (MEK) | 107 | 92 | 49 - 132 | 14 | 19 | | |
| 2-Hexanone | 101 | 89 | 49 - 142 | 12 | 27 | | |
| 4-Methyl-2-pentanone (MIBK) | 104 | 99 | 58 - 136 | 5 | 32 | | |
| Acetone | 84 | 81 | 32 - 126 | 4 | 28 | | |
| Benzene | 94 | 96 | 73 - 121 | 2 | 13 | | |
| Bromoform | 60 | 64 | 45 - 121 | 6 | 19 | | |
| Bromomethane | 54 | 102 | 26 - 136 | 62 | 35 | | F2 |
| Carbon disulfide | 90 | 95 | 54 - 144 | 5 | 34 | | |
| Carbon tetrachloride | 74 | 84 | 65 - 129 | 12 | 20 | | |
| Chlorobenzene | 93 | 94 | 72 - 120 | 0 | 15 | | |
| Dibromochloromethane | 78 | 79 | 62 - 122 | 1 | 19 | | |
| Chloroethane | 62 | 97 | 27 - 131 | 40 | 35 | | F2 |
| Chloroform | 95 | 99 | 73 - 121 | 4 | 17 | | |
| Chloromethane | 95 | 97 | 39 - 134 | 2 | 20 | | |
| cis-1,2-Dichloroethene | 101 | 116 | 66 - 124 | 4 | 22 | | |
| cis-1,3-Dichloropropene | 91 | 92 | 60 - 120 | 1 | 21 | | |
| Dichlorodifluoromethane | 64 | 81 | 14 - 137 | 24 | 34 | | |
| Bromodichloromethane | 85 | 88 | 72 - 120 | 3 | 19 | | |
| Ethylbenzene | 94 | 97 | 68 - 121 | 3 | 16 | | |
| Isopropylbenzene | 94 | 99 | 61 - 122 | 5 | 20 | | |
| Methyl tert-butyl ether | 93 | 95 | 61 - 121 | 2 | 12 | | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212893**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59545-B-2 MS | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC5610.D |
| Dilution: | 500 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1700 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1700 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-59545-B-2 MSD | Analysis Batch: | 240-212893 | Instrument ID: | A3UX15 |
| Client Matrix: | Water | Prep Batch: | N/A | Lab File ID: | UXC5611.D |
| Dilution: | 500 | Leach Batch: | N/A | Initial Weight/Volume: | 5 mL |
| Analysis Date: | 01/04/2016 1723 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 01/04/2016 1723 | | | | 5 mL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Methylene Chloride | 87 | 92 | 70 - 124 | 5 | 14 | | |
| m-Xylene & p-Xylene | 94 | 96 | 66 - 123 | 2 | 15 | | |
| o-Xylene | 95 | 96 | 68 - 121 | 1 | 14 | | |
| Styrene | 96 | 98 | 64 - 126 | 2 | 15 | | |
| Tetrachloroethene | 91 | 92 | 59 - 125 | 2 | 20 | | |
| Toluene | 94 | 92 | 72 - 122 | 2 | 15 | | |
| trans-1,2-Dichloroethene | 98 | 105 | 72 - 125 | 6 | 25 | | |
| trans-1,3-Dichloropropene | 87 | 85 | 58 - 132 | 1 | 22 | | |
| Trichloroethene | 96 | 99 | 61 - 129 | 3 | 14 | | |
| Trichlorofluoromethane | 79 | 95 | 49 - 133 | 18 | 25 | | |
| Vinyl chloride | 95 | 104 | 44 - 122 | 5 | 35 | | |
| Xylenes, Total | 95 | 96 | 67 - 122 | 2 | 14 | | |
| 1,2-Dibromoethane | 101 | 98 | 69 - 125 | 3 | 24 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Toluene-d8 (Surr) | 103 | 98 | 80 - 120 |
| Dibromofluoromethane (Surr) | 100 | 105 | 79 - 120 |
| 4-Bromofluorobenzene (Surr) | 117 | 116 | 61 - 120 |
| 1,2-Dichloroethane-d4 (Surr) | 102 | 97 | 78 - 125 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212893**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59545-B-2 MS | Units: ug/L | MSD Lab Sample ID: | 240-59545-B-2 MSD | |
| Client Matrix: | Water | | Client Matrix: | Water | |
| Dilution: | 500 | | Dilution: | 500 | |
| Analysis Date: | 01/04/2016 1700 | | Analysis Date: | 01/04/2016 1723 | |
| Prep Date: | 01/04/2016 1700 | | Prep Date: | 01/04/2016 1723 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 220 | U | 5000 | 5000 | 4250 | 4480 | |
| 1,1,2,2-Tetrachloroethane | 110 | U | 5000 | 5000 | 4970 | 4970 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 230 | U | 5000 | 5000 | 4400 | 5180 | |
| 1,1-Dichloroethane | 5600 | | 5000 | 5000 | 10600 | 10800 | |
| 1,1-Dichloroethene | 230 | U | 5000 | 5000 | 4740 | 5020 | |
| 1,2,4-Trichlorobenzene | 160 | U | 5000 | 5000 | 3990 | 4280 | |
| 1,2-Dibromo-3-Chloropropane | 410 | U | 5000 | 5000 | 3890 | 3860 | |
| 1,2-Dichlorobenzene | 130 | U | 5000 | 5000 | 4420 | 4610 | |
| 1,2-Dichloroethane | 120 | J | 5000 | 5000 | 4970 | 4960 | |
| 1,2-Dichloropropane | 130 | U | 5000 | 5000 | 4850 | 4910 | |
| 1,3-Dichlorobenzene | 95 | U | 5000 | 5000 | 4440 | 4550 | |
| 1,4-Dichlorobenzene | 140 | U | 5000 | 5000 | 4470 | 4580 | |
| 2-Butanone (MEK) | 660 | J | 10000 | 10000 | 11300 | 9840 | |
| 2-Hexanone | 240 | U | 10000 | 10000 | 10100 | 8950 | |
| 4-Methyl-2-pentanone (MIBK) | 500 | U | 10000 | 10000 | 10400 | 9860 | |
| Acetone | 470 | U | 10000 | 10000 | 8410 | 8100 | |
| Benzene | 180 | U | 5000 | 5000 | 4720 | 4820 | |
| Bromoform | 280 | U | 5000 | 5000 | 3020 | 3200 | |
| Bromomethane | 220 | U | 5000 | 5000 | 2680 | 5100 | F2 |
| Carbon disulfide | 190 | U | 5000 | 5000 | 4490 | 4730 | |
| Carbon tetrachloride | 220 | U | 5000 | 5000 | 3710 | 4200 | |
| Chlorobenzene | 130 | U | 5000 | 5000 | 4670 | 4690 | |
| Dibromochloromethane | 220 | U | 5000 | 5000 | 3910 | 3960 | |
| Chloroethane | 410 | J | 5000 | 5000 | 3500 | 5260 | F2 |
| Chloroform | 130 | U | 5000 | 5000 | 4760 | 4950 | |
| Chloromethane | 220 | U | 5000 | 5000 | 4750 | 4860 | |
| cis-1,2-Dichloroethene | 13000 | | 5000 | 5000 | 17700 | 18400 | |
| cis-1,3-Dichloropropene | 230 | U | 5000 | 5000 | 4550 | 4610 | |
| Dichlorodifluoromethane | 160 | U | 5000 | 5000 | 3190 | 4070 | |
| Bromodichloromethane | 150 | U | 5000 | 5000 | 4260 | 4410 | |
| Ethylbenzene | 130 | U | 5000 | 5000 | 4710 | 4850 | |
| Isopropylbenzene | 180 | U | 5000 | 5000 | 4710 | 4950 | |
| Methyl tert-butyl ether | 100 | U | 5000 | 5000 | 4640 | 4740 | |
| Methylene Chloride | 170 | U | 5000 | 5000 | 4360 | 4590 | |
| m-Xylene & p-Xylene | 120 | U | 5000 | 5000 | 4700 | 4800 | |
| o-Xylene | 130 | U | 5000 | 5000 | 4760 | 4820 | |
| Styrene | 230 | U | 5000 | 5000 | 4790 | 4900 | |
| Tetrachloroethene | 160 | U | 5000 | 5000 | 4540 | 4620 | |
| Toluene | 120 | U | 5000 | 5000 | 4690 | 4620 | |
| trans-1,2-Dichloroethene | 500 | | 5000 | 5000 | 5390 | 5740 | |
| trans-1,3-Dichloropropene | 280 | U | 5000 | 5000 | 4330 | 4270 | |
| Trichloroethene | 110 | U | 5000 | 5000 | 4800 | 4960 | |
| Trichlorofluoromethane | 250 | U | 5000 | 5000 | 3960 | 4740 | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212893**

**Method: 8260C**
**Preparation: 5030C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59545-B-2 MS | Units: | ug/L | MSD Lab Sample ID: | 240-59545-B-2 MSD |
| Client Matrix: | Water | | | Client Matrix: | Water |
| Dilution: | 500 | | | Dilution: | 500 |
| Analysis Date: | 01/04/2016 1700 | | | Analysis Date: | 01/04/2016 1723 |
| Prep Date: | 01/04/2016 1700 | | | Prep Date: | 01/04/2016 1723 |
| Leach Date: | N/A | | | Leach Date: | N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| Vinyl chloride | 4200 | | 5000 | 5000 | 8920 | 9380 |
| Xylenes, Total | 260 | U | 10000 | 10000 | 9460 | 9620 |
| 1,2-Dibromoethane | 160 | U | 5000 | 5000 | 5040 | 4900 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Method Blank - Batch: 240-212548**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-212548/7-A | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-212548 | Lab File ID: | 51231003.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 12/31/2015 1002 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| 1,1'-Biphenyl | 3.5 | U | 3.5 | 50 |
| bis (2-chloroisopropyl) ether | 9.5 | U | 9.5 | 100 |
| 2,4,5-Trichlorophenol | 25 | U | 25 | 150 |
| 2,4,6-Trichlorophenol | 8.9 | U | 8.9 | 150 |
| 2,4-Dichlorophenol | 20 | U | 20 | 150 |
| 2,4-Dimethylphenol | 20 | U | 20 | 150 |
| 2,4-Dinitrophenol | 21 | U | 21 | 330 |
| 2,4-Dinitrotoluene | 17 | U | 17 | 200 |
| 2,6-Dinitrotoluene | 21 | U | 21 | 200 |
| 2-Chloronaphthalene | 0.45 | U | 0.45 | 50 |
| 2-Chlorophenol | 8.2 | U | 8.2 | 50 |
| 2-Methylnaphthalene | 0.50 | U | 0.50 | 6.7 |
| 2-Methylphenol | 11 | U | 11 | 200 |
| 2-Nitroaniline | 9.1 | U | 9.1 | 200 |
| 2-Nitrophenol | 8.3 | U | 8.3 | 50 |
| 3,3'-Dichlorobenzidine | 18 | U | 18 | 100 |
| 3-Nitroaniline | 16 | U | 16 | 200 |
| 4,6-Dinitro-2-methylphenol | 9.2 | U | 9.2 | 150 |
| 4-Bromophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Chloro-3-methylphenol | 21 | U | 21 | 150 |
| 4-Chloroaniline | 17 | U | 17 | 150 |
| 4-Chlorophenyl phenyl ether | 13 | U | 13 | 50 |
| 4-Nitroaniline | 26 | U | 26 | 200 |
| 4-Nitrophenol | 17 | U | 17 | 330 |
| Acenaphthene | 0.76 | U | 0.76 | 6.7 |
| Acenaphthylene | 0.35 | U | 0.35 | 6.7 |
| Acetophenone | 9.2 | U | 9.2 | 100 |
| Anthracene | 0.78 | U | 0.78 | 6.7 |
| Atrazine | 9.1 | U | 9.1 | 200 |
| Benzaldehyde | 12 | U | 12 | 100 |
| Benzo[a]anthracene | 0.63 | U | 0.63 | 6.7 |
| Benzo[a]pyrene | 0.64 | U | 0.64 | 6.7 |
| Benzo[b]fluoranthene | 0.59 | U | 0.59 | 6.7 |
| Benzo[g,h,i]perylene | 0.35 | U | 0.35 | 6.7 |
| Benzo[k]fluoranthene | 0.68 | U | 0.68 | 6.7 |
| Bis(2-chloroethoxy)methane | 22 | U | 22 | 100 |
| Bis(2-chloroethyl)ether | 2.0 | U | 2.0 | 100 |
| Bis(2-ethylhexyl) phthalate | 25.4 | J | 19 | 70 |
| Butyl benzyl phthalate | 10 | U | 10 | 70 |
| Caprolactam | 37 | U | 37 | 330 |
| Carbazole | 27 | U | 27 | 50 |
| Chrysene | 1.1 | U | 1.1 | 6.7 |
| Dibenz(a,h)anthracene | 0.66 | U | 0.66 | 6.7 |
| Dibenzofuran | 0.66 | U | 0.66 | 50 |
| Diethyl phthalate | 16 | U | 16 | 70 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Method Blank - Batch: 240-212548**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-212548/7-A | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-212548 | Lab File ID: | 51231003.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 12/31/2015 1002 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Dimethyl phthalate | 17 | U | 17 | 70 |
| Di-n-butyl phthalate | 15.9 | J | 15 | 70 |
| Di-n-octyl phthalate | 7.9 | U | 7.9 | 70 |
| Fluoranthene | 0.55 | U | 0.55 | 6.7 |
| Fluorene | 0.53 | U | 0.53 | 6.7 |
| Hexachlorobenzene | 2.1 | U | 2.1 | 6.7 |
| Hexachlorobutadiene | 5.6 | U | 5.6 | 50 |
| Hexachlorocyclopentadiene | 8.1 | U | 8.1 | 330 |
| Hexachloroethane | 9.0 | U | 9.0 | 50 |
| Indeno[1,2,3-cd]pyrene | 0.35 | U | 0.35 | 6.7 |
| Isophorone | 13.1 | J | 13 | 50 |
| N-Nitrosodi-n-propylamine | 6.3 | U | 6.3 | 50 |
| N-Nitrosodiphenylamine | 21 | U | 21 | 50 |
| Naphthalene | 0.82 | U | 0.82 | 6.7 |
| Nitrobenzene | 2.2 | U | 2.2 | 100 |
| Pentachlorophenol | 9.1 | U | 9.1 | 150 |
| Phenanthrene | 0.73 | U | 0.73 | 6.7 |
| Phenol | 7.3 | U | 7.3 | 50 |
| Pyrene | 0.44 | U | 0.44 | 6.7 |
| 3 & 4 Methylphenol | 20 | U | 20 | 400 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 89 | 36 - 110 |
| Phenol-d5 (Surr) | 86 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 79 | 20 - 110 |
| 2-Fluorophenol (Surr) | 81 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 78 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 80 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Lab Control Sample - Batch: 240-212548**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-212548/8-A | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-212548 | Lab File ID: | 51231004.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 12/31/2015 1053 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 494 | 74 | 35 - 110 | |
| bis (2-chloroisopropyl) ether | 667 | 493 | 74 | 29 - 110 | |
| 2,4,5-Trichlorophenol | 667 | 502 | 75 | 25 - 110 | |
| 2,4,6-Trichlorophenol | 667 | 520 | 78 | 12 - 110 | |
| 2,4-Dichlorophenol | 667 | 517 | 78 | 39 - 110 | |
| 2,4-Dimethylphenol | 667 | 505 | 76 | 29 - 110 | |
| 2,4-Dinitrophenol | 1330 | 563 | 42 | 10 - 110 | |
| 2,4-Dinitrotoluene | 667 | 582 | 87 | 48 - 110 | |
| 2,6-Dinitrotoluene | 667 | 568 | 85 | 45 - 110 | |
| 2-Chloronaphthalene | 667 | 499 | 75 | 32 - 110 | |
| 2-Chlorophenol | 667 | 524 | 79 | 37 - 110 | |
| 2-Methylnaphthalene | 667 | 514 | 77 | 36 - 110 | |
| 2-Methylphenol | 667 | 526 | 79 | 41 - 110 | |
| 2-Nitroaniline | 667 | 606 | 91 | 45 - 110 | |
| 2-Nitrophenol | 667 | 577 | 87 | 34 - 110 | |
| 3,3'-Dichlorobenzidine | 1330 | 1100 | 82 | 28 - 110 | |
| 3-Nitroaniline | 667 | 536 | 80 | 44 - 110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 1000 | 75 | 10 - 110 | |
| 4-Bromophenyl phenyl ether | 667 | 554 | 83 | 39 - 110 | |
| 4-Chloro-3-methylphenol | 667 | 558 | 84 | 48 - 110 | |
| 4-Chloroaniline | 667 | 441 | 66 | 30 - 110 | |
| 4-Chlorophenyl phenyl ether | 667 | 513 | 77 | 40 - 110 | |
| 4-Nitroaniline | 667 | 630 | 94 | 48 - 110 | |
| 4-Nitrophenol | 1330 | 1450 | 109 | 28 - 110 | |
| Acenaphthene | 667 | 501 | 75 | 38 - 110 | |
| Acenaphthylene | 667 | 491 | 74 | 40 - 110 | |
| Acetophenone | 667 | 498 | 75 | 40 - 110 | |
| Anthracene | 667 | 528 | 79 | 48 - 110 | |
| Atrazine | 1330 | 1330 | 100 | 54 - 120 | |
| Benzaldehyde | 1330 | 1090 | 82 | 32 - 110 | |
| Benzo[a]anthracene | 667 | 537 | 80 | 50 - 110 | |
| Benzo[a]pyrene | 667 | 522 | 78 | 44 - 110 | |
| Benzo[b]fluoranthene | 667 | 524 | 79 | 43 - 110 | |
| Benzo[g,h,i]perylene | 667 | 567 | 85 | 51 - 110 | |
| Benzo[k]fluoranthene | 667 | 525 | 79 | 38 - 105 | |
| Bis(2-chloroethoxy)methane | 667 | 546 | 82 | 32 - 110 | |
| Bis(2-chloroethyl)ether | 667 | 501 | 75 | 34 - 110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 642 | 96 | 50 - 110 | |
| Butyl benzyl phthalate | 667 | 609 | 91 | 51 - 110 | |
| Caprolactam | 1330 | 1140 | 85 | 44 - 114 | |
| Carbazole | 667 | 638 | 96 | 50 - 110 | |
| Chrysene | 667 | 544 | 82 | 50 - 110 | |
| Dibenz(a,h)anthracene | 667 | 581 | 87 | 51 - 110 | |
| Dibenzofuran | 667 | 503 | 75 | 43 - 110 | |
| Diethyl phthalate | 667 | 563 | 84 | 52 - 110 | |
| Dimethyl phthalate | 667 | 550 | 82 | 50 - 110 | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Lab Control Sample - Batch: 240-212548**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-212548/8-A | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-212548 | Lab File ID: | 51231004.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 12/31/2015 1053 | Units: | ug/Kg | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Di-n-butyl phthalate | 667 | 708 | 106 | 51 - 110 | |
| Di-n-octyl phthalate | 667 | 571 | 86 | 48 - 110 | |
| Fluoranthene | 667 | 589 | 88 | 51 - 110 | |
| Fluorene | 667 | 465 | 70 | 46 - 110 | |
| Hexachlorobenzene | 667 | 523 | 79 | 43 - 110 | |
| Hexachlorobutadiene | 667 | 515 | 77 | 29 - 110 | |
| Hexachlorocyclopentadiene | 667 | 399 | 60 | 12 - 110 | |
| Hexachloroethane | 667 | 520 | 78 | 30 - 110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 574 | 86 | 50 - 110 | |
| Isophorone | 667 | 555 | 83 | 36 - 110 | |
| N-Nitrosodi-n-propylamine | 667 | 544 | 82 | 38 - 110 | |
| N-Nitrosodiphenylamine | 1330 | 1100 | 82 | 46 - 110 | |
| Naphthalene | 667 | 501 | 75 | 36 - 110 | |
| Nitrobenzene | 667 | 545 | 82 | 32 - 110 | |
| Pentachlorophenol | 1330 | 1110 | 83 | 10 - 110 | |
| Phenanthrene | 667 | 530 | 79 | 49 - 110 | |
| Phenol | 667 | 505 | 76 | 38 - 110 | |
| Pyrene | 667 | 574 | 86 | 49 - 110 | |
| 3 & 4 Methylphenol | 667 | 516 | 77 | 40 - 110 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Terphenyl-d14 (Surr) | 81 | 36 - 110 |
| Phenol-d5 (Surr) | 80 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 82 | 20 - 110 |
| 2-Fluorophenol (Surr) | 77 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 74 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 90 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212548**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-212548 | Lab File ID: | 51231009.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.67 g |
| Analysis Date: | 12/31/2015 1258 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-212548 | Lab File ID: | 51231010.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30.03 g |
| Analysis Date: | 12/31/2015 1323 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| 1,1'-Biphenyl | 76 | 78 | 32 - 110 | 1 | 32 | | |
| bis (2-chloroisopropyl) ether | 73 | 76 | 11 - 110 | 2 | 42 | | |
| 2,4,5-Trichlorophenol | 76 | 79 | 10 - 117 | 3 | 99 | | |
| 2,4,6-Trichlorophenol | 81 | 78 | 10 - 110 | 5 | 38 | | |
| 2,4-Dichlorophenol | 77 | 79 | 10 - 110 | 1 | 34 | | |
| 2,4-Dimethylphenol | 87 | 84 | 10 - 110 | 5 | 31 | | |
| 2,4-Dinitrophenol | 94 | 89 | 10 - 110 | 7 | 99 | | |
| 2,4-Dinitrotoluene | 95 | 98 | 32 - 110 | 2 | 30 | | |
| 2,6-Dinitrotoluene | 91 | 91 | 35 - 110 | 1 | 30 | | |
| 2-Chloronaphthalene | 77 | 78 | 28 - 110 | 0 | 30 | | |
| 2-Chlorophenol | 76 | 78 | 10 - 110 | 1 | 47 | | |
| 2-Methylnaphthalene | 76 | 78 | 10 - 133 | 2 | 42 | | |
| 2-Methylphenol | 75 | 78 | 24 - 110 | 2 | 51 | | |
| 2-Nitroaniline | 98 | 97 | 39 - 110 | 2 | 31 | | |
| 2-Nitrophenol | 86 | 92 | 10 - 110 | 6 | 49 | | |
| 3,3'-Dichlorobenzidine | 80 | 90 | 10 - 110 | 11 | 56 | | |
| 3-Nitroaniline | 59 | 72 | 10 - 110 | 18 | 30 | | |
| 4,6-Dinitro-2-methylphenol | 98 | 97 | 10 - 110 | 2 | 55 | | |
| 4-Bromophenyl phenyl ether | 93 | 90 | 33 - 110 | 4 | 30 | | |
| 4-Chloro-3-methylphenol | 88 | 89 | 25 - 110 | 0 | 54 | | |
| 4-Chloroaniline | 48 | 46 | 10 - 110 | 5 | 36 | | |
| 4-Chlorophenyl phenyl ether | 81 | 81 | 32 - 110 | 1 | 30 | | |
| 4-Nitroaniline | 80 | 98 | 10 - 110 | 18 | 48 | | |
| 4-Nitrophenol | 122 | 107 | 10 - 113 | 14 | 49 | F1 | |
| Acenaphthene | 78 | 78 | 22 - 110 | 2 | 99 | | |
| Acenaphthylene | 79 | 79 | 24 - 110 | 1 | 99 | | |
| Acetophenone | 79 | 79 | 31 - 110 | 1 | 43 | | |
| Anthracene | 86 | 86 | 20 - 110 | 1 | 99 | | |
| Atrazine | 105 | 107 | 45 - 118 | 1 | 30 | | |
| Benzaldehyde | 81 | 78 | 23 - 110 | 4 | 42 | | |
| Benzo[a]anthracene | 86 | 90 | 10 - 122 | 3 | 99 | | |
| Benzo[a]pyrene | 86 | 87 | 10 - 110 | 0 | 99 | | |
| Benzo[b]fluoranthene | 90 | 90 | 12 - 118 | 1 | 99 | | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212548**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-212548 | Lab File ID: | 51231009.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.67 g |
| Analysis Date: | 12/31/2015 1258 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212723 | Instrument ID: | A4HP9 |
| Client Matrix: | Solid | Prep Batch: | 240-212548 | Lab File ID: | 51231010.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30.03 g |
| Analysis Date: | 12/31/2015 1323 | | | Final Weight/Volume: | 2 mL |
| Prep Date: | 12/30/2015 0705 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | | |

| Analyte | % Rec. MS | % Rec. MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
|---|---|---|---|---|---|---|---|
| Benzo[g,h,i]perylene | 99 | 96 | 10 - 117 | 4 | 99 | | |
| Benzo[k]fluoranthene | 86 | 88 | 10 - 121 | 1 | 99 | | |
| Bis(2-chloroethoxy)methane | 82 | 84 | 26 - 110 | 0 | 37 | | |
| Bis(2-chloroethyl)ether | 84 | 78 | 21 - 110 | 9 | 55 | | |
| Bis(2-ethylhexyl) phthalate | 85 | 95 | 40 - 110 | 9 | 30 | | |
| Butyl benzyl phthalate | 100 | 102 | 44 - 110 | 1 | 30 | | |
| Caprolactam | 90 | 92 | 10 - 134 | 1 | 32 | | |
| Carbazole | 92 | 101 | 34 - 110 | 8 | 30 | | |
| Chrysene | 87 | 90 | 10 - 125 | 2 | 99 | | |
| Dibenz(a,h)anthracene | 102 | 99 | 14 - 113 | 5 | 99 | | |
| Dibenzofuran | 81 | 80 | 29 - 110 | 2 | 30 | | |
| Diethyl phthalate | 94 | 94 | 42 - 110 | 2 | 30 | | |
| Dimethyl phthalate | 90 | 89 | 41 - 110 | 1 | 30 | | |
| Di-n-butyl phthalate | 110 | 111 | 43 - 110 | 0 | 30 | | F1 |
| Di-n-octyl phthalate | 96 | 97 | 24 - 119 | 1 | 30 | | |
| Fluoranthene | 93 | 95 | 10 - 110 | 1 | 99 | | |
| Fluorene | 76 | 76 | 23 - 110 | 2 | 99 | | |
| Hexachlorobenzene | 86 | 83 | 34 - 110 | 4 | 30 | | |
| Hexachlorobutadiene | 76 | 84 | 25 - 110 | 8 | 34 | | |
| Hexachlorocyclopentadiene | 40 | 37 | 10 - 110 | 10 | 79 | J | J |
| Hexachloroethane | 75 | 79 | 12 - 110 | 5 | 50 | | |
| Indeno[1,2,3-cd]pyrene | 98 | 96 | 10 - 114 | 4 | 99 | | |
| Isophorone | 79 | 82 | 29 - 110 | 2 | 38 | | |
| N-Nitrosodi-n-propylamine | 82 | 84 | 26 - 110 | 1 | 42 | | |
| N-Nitrosodiphenylamine | 87 | 87 | 22 - 110 | 2 | 30 | | |
| Naphthalene | 74 | 76 | 10 - 111 | 2 | 99 | | |
| Nitrobenzene | 82 | 83 | 23 - 110 | 1 | 41 | | |
| Pentachlorophenol | 92 | 79 | 10 - 110 | 17 | 50 | | |
| Phenanthrene | 85 | 85 | 10 - 166 | 1 | 99 | | |
| Phenol | 73 | 76 | 17 - 110 | 4 | 53 | | |
| Pyrene | 93 | 93 | 10 - 147 | 1 | 99 | | |
| 3 & 4 Methylphenol | 74 | 76 | 25 - 110 | 1 | 50 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|

Client: Tetra Tech, Inc. Job Number: 240-59482-1

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Terphenyl-d14 (Surr) | 89 | 89 | 36 - 110 |
| Phenol-d5 (Surr) | 80 | 79 | 26 - 110 |
| Nitrobenzene-d5 (Surr) | 84 | 86 | 20 - 110 |
| 2-Fluorophenol (Surr) | 78 | 77 | 24 - 110 |
| 2-Fluorobiphenyl (Surr) | 77 | 78 | 24 - 110 |
| 2,4,6-Tribromophenol (Surr) | 101 | 90 | 10 - 110 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212548**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Units: ug/Kg | MSD Lab Sample ID: | 240-59482-2 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 12/31/2015 1258 | | Analysis Date: | 12/31/2015 1323 | |
| Prep Date: | 12/30/2015 0705 | | Prep Date: | 12/30/2015 0705 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual |
|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 4.3 | U | 826 | 816 | 631 | | 636 |
| bis (2-chloroisopropyl) ether | 12 | U | 826 | 816 | 607 | | 617 |
| 2,4,5-Trichlorophenol | 31 | U | 826 | 816 | 627 | | 643 |
| 2,4,6-Trichlorophenol | 11 | U | 826 | 816 | 670 | | 636 |
| 2,4-Dichlorophenol | 24 | U | 826 | 816 | 640 | | 646 |
| 2,4-Dimethylphenol | 24 | U | 826 | 816 | 717 | | 684 |
| 2,4-Dinitrophenol | 26 | U | 1650 | 1630 | 1550 | | 1450 |
| 2,4-Dinitrotoluene | 21 | U | 826 | 816 | 788 | | 801 |
| 2,6-Dinitrotoluene | 26 | U | 826 | 816 | 752 | | 742 |
| 2-Chloronaphthalene | 0.55 | U | 826 | 816 | 636 | | 636 |
| 2-Chlorophenol | 10 | U | 826 | 816 | 631 | | 640 |
| 2-Methylnaphthalene | 0.61 | U | 826 | 816 | 629 | | 641 |
| 2-Methylphenol | 13 | U | 826 | 816 | 623 | | 638 |
| 2-Nitroaniline | 11 | U | 826 | 816 | 808 | | 791 |
| 2-Nitrophenol | 10 | U | 826 | 816 | 710 | | 752 |
| 3,3'-Dichlorobenzidine | 22 | U | 1650 | 1630 | 1320 | | 1470 |
| 3-Nitroaniline | 20 | U | 826 | 816 | 490 | | 589 |
| 4,6-Dinitro-2-methylphenol | 11 | U | 1650 | 1630 | 1620 | | 1590 |
| 4-Bromophenyl phenyl ether | 16 | U | 826 | 816 | 767 | | 736 |
| 4-Chloro-3-methylphenol | 26 | U | 826 | 816 | 727 | | 726 |
| 4-Chloroaniline | 21 | U | 826 | 816 | 396 | | 378 |
| 4-Chlorophenyl phenyl ether | 16 | U | 826 | 816 | 672 | | 665 |
| 4-Nitroaniline | 32 | U | 826 | 816 | 664 | | 799 |
| 4-Nitrophenol | 21 | U | 1650 | 1630 | 2010 | F1 | 1740 |
| Acenaphthene | 0.93 | U | 826 | 816 | 648 | | 636 |
| Acenaphthylene | 0.43 | U | 826 | 816 | 654 | | 646 |
| Acetophenone | 11 | U | 826 | 816 | 652 | | 648 |
| Anthracene | 0.95 | U | 826 | 816 | 711 | | 704 |
| Atrazine | 11 | U | 1650 | 1630 | 1730 | | 1740 |
| Benzaldehyde | 15 | U | 1650 | 1630 | 1330 | | 1270 |
| Benzo[a]anthracene | 0.77 | U | 826 | 816 | 711 | | 731 |
| Benzo[a]pyrene | 0.78 | U | 826 | 816 | 713 | | 714 |
| Benzo[b]fluoranthene | 0.72 | U | 826 | 816 | 746 | | 738 |
| Benzo[g,h,i]perylene | 0.43 | U | 826 | 816 | 819 | | 783 |
| Benzo[k]fluoranthene | 0.83 | U | 826 | 816 | 712 | | 716 |
| Bis(2-chloroethoxy)methane | 27 | U | 826 | 816 | 681 | | 683 |
| Bis(2-chloroethyl)ether | 2.4 | U | 826 | 816 | 697 | | 637 |
| Bis(2-ethylhexyl) phthalate | 160 | | 826 | 816 | 861 | | 938 |
| Butyl benzyl phthalate | 12 | U | 826 | 816 | 822 | | 832 |
| Caprolactam | 45 | U | 1650 | 1630 | 1490 | | 1500 |
| Carbazole | 33 | U | 826 | 816 | 763 | | 828 |
| Chrysene | 1.3 | U | 826 | 816 | 717 | | 734 |
| Dibenz(a,h)anthracene | 0.81 | U | 826 | 816 | 844 | | 805 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212548**

**Method: 8270D**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Units: ug/Kg | MSD Lab Sample ID: | 240-59482-2 | |
| Client Matrix: | Solid | | Client Matrix: | Solid | |
| Dilution: | 1.0 | | Dilution: | 1.0 | |
| Analysis Date: | 12/31/2015 1258 | | Analysis Date: | 12/31/2015 1323 | |
| Prep Date: | 12/30/2015 0705 | | Prep Date: | 12/30/2015 0705 | |
| Leach Date: | N/A | | Leach Date: | N/A | |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|---|
| Dibenzofuran | 0.81 | U | 826 | 816 | 669 | | 656 | |
| Diethyl phthalate | 20 | U | 826 | 816 | 779 | | 764 | |
| Dimethyl phthalate | 21 | U | 826 | 816 | 741 | | 730 | |
| Di-n-butyl phthalate | 30 | J | 826 | 816 | 937 | | 939 | F1 |
| Di-n-octyl phthalate | 9.7 | U | 826 | 816 | 794 | | 788 | |
| Fluoranthene | 9.3 | | 826 | 816 | 781 | | 787 | |
| Fluorene | 0.65 | U | 826 | 816 | 631 | | 619 | |
| Hexachlorobenzene | 2.6 | U | 826 | 816 | 709 | | 682 | |
| Hexachlorobutadiene | 6.9 | U | 826 | 816 | 629 | | 683 | |
| Hexachlorocyclopentadiene | 9.9 | U | 826 | 816 | 333 | J | 302 | J |
| Hexachloroethane | 11 | U | 826 | 816 | 617 | | 646 | |
| Indeno[1,2,3-cd]pyrene | 0.43 | U | 826 | 816 | 811 | | 782 | |
| Isophorone | 21 | J | 826 | 816 | 676 | | 691 | |
| N-Nitrosodi-n-propylamine | 7.7 | U | 826 | 816 | 674 | | 684 | |
| N-Nitrosodiphenylamine | 26 | U | 1650 | 1630 | 1440 | | 1420 | |
| Naphthalene | 4.3 | J | 826 | 816 | 612 | | 623 | |
| Nitrobenzene | 2.7 | U | 826 | 816 | 676 | | 680 | |
| Pentachlorophenol | 11 | U | 1650 | 1630 | 1520 | | 1280 | |
| Phenanthrene | 9.3 | | 826 | 816 | 715 | | 706 | |
| Phenol | 8.9 | U | 826 | 816 | 601 | | 623 | |
| Pyrene | 7.9 | J | 826 | 816 | 773 | | 766 | |
| 3 & 4 Methylphenol | 24 | U | 826 | 816 | 613 | | 618 | |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Method Blank - Batch: 240-212559**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-212559/1-A | Analysis Batch: | 240-212533 | Instrument ID: | YPID |
| Client Matrix: | Solid | Prep Batch: | 240-212559 | Lab File ID: | Y0122919.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 12/30/2015 0918 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 0739 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 46 | U | 46 | 100 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 68 | 40 - 139 |

**Lab Control Sample - Batch: 240-212559**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-212559/2-A | Analysis Batch: | 240-212533 | Instrument ID: | YPID |
| Client Matrix: | Solid | Prep Batch: | 240-212559 | Lab File ID: | Y0122920.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5 g |
| Analysis Date: | 12/30/2015 0950 | Units: | ug/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 0739 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 771 | 96 | 74 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Trifluorotoluene (Surr) | 90 | 40 - 139 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212559**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212533 | Instrument ID: | YPID |
| Client Matrix: | Solid | Prep Batch: | 240-212559 | Lab File ID: | Y0122922.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.01 g |
| Analysis Date: | 12/30/2015 1057 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 0739 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-59482-2 | Analysis Batch: | 240-212533 | Instrument ID: | YPID |
| Client Matrix: | Solid | Prep Batch: | 240-212559 | Lab File ID: | Y0122923.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 5.04 g |
| Analysis Date: | 12/30/2015 1131 | | | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 0739 | | | Injection Volume: | 5 mL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | % Rec. | | | | | | |
|---|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Gasoline Range Organics [C6 - C10] | 90 | 86 | 10 - 136 | 5 | 40 | | |

| Surrogate | MS % Rec | MSD % Rec | Acceptance Limits |
|---|---|---|---|
| Trifluorotoluene (Surr) | 82 | 82 | 40 - 139 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212559**

**Method: 8015C**
**Preparation: 5030A**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59482-2 | Units: | ug/Kg | MSD Lab Sample ID: | 240-59482-2 |
| Client Matrix: | Solid | | | Client Matrix: | Solid |
| Dilution: | 1.0 | | | Dilution: | 1.0 |
| Analysis Date: | 12/30/2015 1057 | | | Analysis Date: | 12/30/2015 1131 |
| Prep Date: | 12/30/2015 0739 | | | Prep Date: | 12/30/2015 0739 |
| Leach Date: | N/A | | | Leach Date: | N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual |
|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 56 | U | 979 | 973 | 885 | 841 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Method Blank - Batch: 240-212549**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-212549/8-A | Analysis Batch: | 240-212780 | Instrument ID: | A2HP5F |
| Client Matrix: | Solid | Prep Batch: | 240-212549 | Lab File ID: | P5F23115.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 12/31/2015 2019 | Units: | mg/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 0714 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Total Petroleum Hydrocarbons (C10-C32) | 6.4 | U | 6.4 | 17 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| o-Terphenyl (Surr) | 90 | 40 - 160 |

**Lab Control Sample - Batch: 240-212549**

**Method: 8015C**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-212549/9-A | Analysis Batch: | 240-212780 | Instrument ID: | A2HP5F |
| Client Matrix: | Solid | Prep Batch: | 240-212549 | Lab File ID: | P5F23116.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 12/31/2015 2050 | Units: | mg/Kg | Final Weight/Volume: | 5 mL |
| Prep Date: | 12/30/2015 0714 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Diesel | 84.0 | 75.4 | 90 | 52 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| o-Terphenyl (Surr) | 97 | 40 - 160 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Method Blank - Batch: 240-212544**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | MB 240-212544/7-A | Analysis Batch: | 240-212845 | Instrument ID: | A2HP12 |
| Client Matrix: | Solid | Prep Batch: | 240-212544 | Lab File ID: | P1200019.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 01/04/2016 1243 | Units: | ug/Kg | Final Weight/Volume: | 10 mL |
| Prep Date: | 12/30/2015 0654 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Result | Qual | MDL | RL |
|---|---|---|---|---|
| Aroclor-1016 | 12 | U | 12 | 33 |
| Aroclor-1221 | 16 | U | 16 | 33 |
| Aroclor-1232 | 20 | U | 20 | 33 |
| Aroclor-1242 | 11 | U | 11 | 33 |
| Aroclor-1248 | 8.0 | U | 8.0 | 33 |
| Aroclor-1254 | 14 | U | 14 | 33 |
| Aroclor-1260 | 9.0 | U | 9.0 | 33 |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Tetrachloro-m-xylene | 106 | 13 - 134 |
| DCB Decachlorobiphenyl | 110 | 10 - 155 |

**Lab Control Sample - Batch: 240-212544**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | LCS 240-212544/8-A | Analysis Batch: | 240-212845 | Instrument ID: | A2HP12 |
| Client Matrix: | Solid | Prep Batch: | 240-212544 | Lab File ID: | P1200020.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 30 g |
| Analysis Date: | 01/04/2016 1301 | Units: | ug/Kg | Final Weight/Volume: | 10 mL |
| Prep Date: | 12/30/2015 0654 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| Analyte | Spike Amount | Result | % Rec. | Limit | Qual |
|---|---|---|---|---|---|
| Aroclor-1016 | 333 | 314 | 94 | 51 - 120 | |
| Aroclor-1260 | 333 | 337 | 101 | 48 - 120 | |

| Surrogate | % Rec | Acceptance Limits |
|---|---|---|
| Tetrachloro-m-xylene | 102 | 13 - 134 |
| DCB Decachlorobiphenyl | 110 | 10 - 155 |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212544**

**Method: 8082A**
**Preparation: 3540C**

| | | | | | |
|---|---|---|---|---|---|
| MS Lab Sample ID: | 240-59503-G-4-B MS | Analysis Batch: | 240-212845 | Instrument ID: | A2HP12 |
| Client Matrix: | Solid | Prep Batch: | 240-212544 | Lab File ID: | P1200017.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.79 g |
| Analysis Date: | 01/04/2016 1207 | | | Final Weight/Volume: | 10 mL |
| Prep Date: | 12/30/2015 0654 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | | | | | |
|---|---|---|---|---|---|
| MSD Lab Sample ID: | 240-59503-G-4-C MSD | Analysis Batch: | 240-212845 | Instrument ID: | A2HP12 |
| Client Matrix: | Solid | Prep Batch: | 240-212544 | Lab File ID: | P1200018.D |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | 29.86 g |
| Analysis Date: | 01/04/2016 1225 | | | Final Weight/Volume: | 10 mL |
| Prep Date: | 12/30/2015 0654 | | | Injection Volume: | 1 uL |
| Leach Date: | N/A | | | Column ID: | PRIMARY |

| | % Rec. | | | | | |
|---|---|---|---|---|---|---|
| Analyte | MS | MSD | Limit | RPD | RPD Limit | MS Qual | MSD Qual |
| Aroclor-1016 | 40 | 33 | 17 - 141 | 17 | 56 | | |
| Aroclor-1260 | 32 | 27 | 19 - 137 | 29 | 55 | | p |

| Surrogate | MS % Rec | MSD % Rec | | Acceptance Limits |
|---|---|---|---|---|
| Tetrachloro-m-xylene | 44 | 31 | p | 13 - 134 |
| DCB Decachlorobiphenyl | 43 | 37 | | 10 - 155 |

**Matrix Spike/**
**Matrix Spike Duplicate Recovery Report - Batch: 240-212544**

**Method: 8082A**
**Preparation: 3540C**

| | | | |
|---|---|---|---|
| MS Lab Sample ID: | 240-59503-G-4-B MS | Units: ug/Kg | MSD Lab Sample ID: 240-59503-G-4-C MSD |
| Client Matrix: | Solid | | Client Matrix: Solid |
| Dilution: | 1.0 | | Dilution: 1.0 |
| Analysis Date: | 01/04/2016 1207 | | Analysis Date: 01/04/2016 1225 |
| Prep Date: | 12/30/2015 0654 | | Prep Date: 12/30/2015 0654 |
| Leach Date: | N/A | | Leach Date: N/A |

| Analyte | Sample Result/Qual | | MS Spike Amount | MSD Spike Amount | MS Result/Qual | MSD Result/Qual | |
|---|---|---|---|---|---|---|---|
| Aroclor-1016 | 16 | U | 442 | 441 | 175 | 148 | |
| Aroclor-1260 | 12 | U | 442 | 441 | 143 | 118 | p |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

**Duplicate - Batch: 240-212487**

**Method: Moisture**
**Preparation: N/A**

| | | | | | |
|---|---|---|---|---|---|
| Lab Sample ID: | 240-59255-B-13 DU | Analysis Batch: | 240-212487 | Instrument ID: | No Equipment Assigned |
| Client Matrix: | Solid | Prep Batch: | N/A | Lab File ID: | N/A |
| Dilution: | 1.0 | Leach Batch: | N/A | Initial Weight/Volume: | |
| Analysis Date: | 12/29/2015 1350 | Units: | % | Final Weight/Volume: | |
| Prep Date: | N/A | | | | |
| Leach Date: | N/A | | | | |

| Analyte | Sample Result/Qual | Result | RPD | Limit | Qual |
|---|---|---|---|---|---|
| Percent Solids | 98 | 98 | 0.08 | 20 | |
| Percent Moisture | 1.7 | 1.7 | 5 | 20 | |

# DATA REPORTING QUALIFIERS

Client: Tetra Tech, Inc. Job Number: 240-59482-1

| Lab Section | Qualifier | Description |
|---|---|---|
| GC/MS VOA | | |
| | B | Compound was found in the blank and sample. |
| | U | Indicates the analyte was analyzed for but not detected. |
| | F1 | MS and/or MSD Recovery is outside acceptance limits. |
| | F2 | MS/MSD RPD exceeds control limits |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| GC/MS Semi VOA | | |
| | B | Compound was found in the blank and sample. |
| | U | Indicates the analyte was analyzed for but not detected. |
| | F1 | MS and/or MSD Recovery is outside acceptance limits. |
| | J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| GC VOA | | |
| | U | Indicates the analyte was analyzed for but not detected. |
| GC Semi VOA | | |
| | U | Indicates the analyte was analyzed for but not detected. |
| | p | The %RPD between the primary and confirmation column/detector is >40%. The lower value has been reported. |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC/MS VOA** | | | | | |
| **Analysis Batch:240-212606** | | | | | |
| LCS 240-212606/23 | Lab Control Sample | T | Solid | 8260C | |
| LCS 240-212606/25 | Lab Control Sample | T | Solid | 8260C | |
| MB 240-212606/24 | Method Blank | T | Solid | 8260C | |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 8260C | 240-212657 |
| 240-59482-2MS | Matrix Spike | T | Solid | 8260C | 240-212657 |
| 240-59482-2MSD | Matrix Spike Duplicate | T | Solid | 8260C | 240-212657 |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 8260C | 240-212657 |
| **Prep Batch: 240-212657** | | | | | |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 5030C | |
| 240-59482-2MS | Matrix Spike | T | Solid | 5030C | |
| 240-59482-2MSD | Matrix Spike Duplicate | T | Solid | 5030C | |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 5030C | |
| **Analysis Batch:240-212893** | | | | | |
| LCS 240-212893/5 | Lab Control Sample | T | Water | 8260C | |
| MB 240-212893/8 | Method Blank | T | Water | 8260C | |
| 240-59482-1 | TB-122815 | T | Water | 8260C | |
| 240-59545-B-2 MS | Matrix Spike | T | Water | 8260C | |
| 240-59545-B-2 MSD | Matrix Spike Duplicate | T | Water | 8260C | |

**Report Basis**
T = Total

| | | | | | |
|---|---|---|---|---|---|
| **GC/MS Semi VOA** | | | | | |
| **Prep Batch: 240-212548** | | | | | |
| LCS 240-212548/8-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-212548/7-A | Method Blank | T | Solid | 3540C | |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 3540C | |
| 240-59482-2MS | Matrix Spike | T | Solid | 3540C | |
| 240-59482-2MSD | Matrix Spike Duplicate | T | Solid | 3540C | |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 3540C | |
| **Analysis Batch:240-212723** | | | | | |
| LCS 240-212548/8-A | Lab Control Sample | T | Solid | 8270D | 240-212548 |
| MB 240-212548/7-A | Method Blank | T | Solid | 8270D | 240-212548 |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 8270D | 240-212548 |
| 240-59482-2MS | Matrix Spike | T | Solid | 8270D | 240-212548 |
| 240-59482-2MSD | Matrix Spike Duplicate | T | Solid | 8270D | 240-212548 |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 8270D | 240-212548 |

**Report Basis**
T = Total

**TestAmerica Canton**

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC VOA** | | | | | |
| **Analysis Batch:240-212533** | | | | | |
| LCS 240-212559/2-A | Lab Control Sample | T | Solid | 8015C | 240-212559 |
| MB 240-212559/1-A | Method Blank | T | Solid | 8015C | 240-212559 |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 8015C | 240-212559 |
| 240-59482-2MS | Matrix Spike | T | Solid | 8015C | 240-212559 |
| 240-59482-2MSD | Matrix Spike Duplicate | T | Solid | 8015C | 240-212559 |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 8015C | 240-212559 |
| **Prep Batch: 240-212559** | | | | | |
| LCS 240-212559/2-A | Lab Control Sample | T | Solid | 5030A | |
| MB 240-212559/1-A | Method Blank | T | Solid | 5030A | |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 5030A | |
| 240-59482-2MS | Matrix Spike | T | Solid | 5030A | |
| 240-59482-2MSD | Matrix Spike Duplicate | T | Solid | 5030A | |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 5030A | |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## QC Association Summary

| Lab Sample ID | Client Sample ID | Report Basis | Client Matrix | Method | Prep Batch |
|---|---|---|---|---|---|
| **GC Semi VOA** | | | | | |
| **Prep Batch: 240-212544** | | | | | |
| LCS 240-212544/8-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-212544/7-A | Method Blank | T | Solid | 3540C | |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 3540C | |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 3540C | |
| 240-59503-G-4-B MS | Matrix Spike | T | Solid | 3540C | |
| 240-59503-G-4-C MSD | Matrix Spike Duplicate | T | Solid | 3540C | |
| **Prep Batch: 240-212549** | | | | | |
| LCS 240-212549/9-A | Lab Control Sample | T | Solid | 3540C | |
| MB 240-212549/8-A | Method Blank | T | Solid | 3540C | |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 3540C | |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 3540C | |
| **Analysis Batch:240-212780** | | | | | |
| LCS 240-212549/9-A | Lab Control Sample | T | Solid | 8015C | 240-212549 |
| MB 240-212549/8-A | Method Blank | T | Solid | 8015C | 240-212549 |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 8015C | 240-212549 |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 8015C | 240-212549 |
| **Analysis Batch:240-212845** | | | | | |
| LCS 240-212544/8-A | Lab Control Sample | T | Solid | 8082A | 240-212544 |
| MB 240-212544/7-A | Method Blank | T | Solid | 8082A | 240-212544 |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | 8082A | 240-212544 |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | 8082A | 240-212544 |
| 240-59503-G-4-B MS | Matrix Spike | T | Solid | 8082A | 240-212544 |
| 240-59503-G-4-C MSD | Matrix Spike Duplicate | T | Solid | 8082A | 240-212544 |

**Report Basis**
T = Total

**General Chemistry**

| | | | | | |
|---|---|---|---|---|---|
| **Analysis Batch:240-212487** | | | | | |
| 240-59255-B-13 DU | Duplicate | T | Solid | Moisture | |
| 240-59482-2 | F-VB-UST-P-7 | T | Solid | Moisture | |
| 240-59482-3 | F-VB-UST-P-8 | T | Solid | Moisture | |

**Report Basis**
T = Total

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Laboratory Chronicle

**Lab ID: 240-59482-1** **Client ID: TB-122815**

Sample Date/Time: 12/28/2015 00:00 Received Date/Time: 12/29/2015 09:30

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-59482-B-1 | | 240-212893 | | 01/04/2016 12:07 | 1 | TAL CAN | LEE |
| A:8260C | 240-59482-B-1 | | 240-212893 | | 01/04/2016 12:07 | 1 | TAL CAN | LEE |

**Lab ID: 240-59482-2** **Client ID: F-VB-UST-P-7**

Sample Date/Time: 12/28/2015 10:00 Received Date/Time: 12/29/2015 09:30

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-59482-B-2-A | | 240-212606 | 240-212657 | 12/30/2015 14:25 | 1 | TAL CAN | SAM |
| A:8260C | 240-59482-B-2-A | | 240-212606 | 240-212657 | 12/30/2015 21:48 | 1 | TAL CAN | SAM |
| P:3540C | 240-59482-C-2-B | | 240-212723 | 240-212548 | 12/30/2015 07:05 | 1 | TAL CAN | YDN |
| A:8270D | 240-59482-C-2-B | | 240-212723 | 240-212548 | 12/31/2015 12:33 | 1 | TAL CAN | MRU |
| P:5030A | 240-59482-C-2-F | | 240-212533 | 240-212559 | 12/30/2015 07:39 | 1 | TAL CAN | RTR |
| A:8015C | 240-59482-C-2-F | | 240-212533 | 240-212559 | 12/30/2015 10:23 | 1 | TAL CAN | RTR |
| P:3540C | 240-59482-C-2-E | | 240-212780 | 240-212549 | 12/30/2015 07:14 | 1 | TAL CAN | YDN |
| A:8015C | 240-59482-C-2-E | | 240-212780 | 240-212549 | 12/31/2015 17:44 | 1 | TAL CAN | HMB |
| P:3540C | 240-59482-C-2-A | | 240-212845 | 240-212544 | 12/30/2015 06:54 | 1 | TAL CAN | YDN |
| A:8082A | 240-59482-C-2-A | | 240-212845 | 240-212544 | 01/04/2016 10:19 | 1 | TAL CAN | HMB |
| A:Moisture | 240-59482-A-2 | | 240-212487 | | 12/29/2015 13:50 | 1 | TAL CAN | BLW |

**Lab ID: 240-59482-2 MS** **Client ID: F-VB-UST-P-7**

Sample Date/Time: 12/28/2015 10:00 Received Date/Time: 12/29/2015 09:30

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-59482-B-2-B MS | | 240-212606 | 240-212657 | 12/30/2015 14:25 | 1 | TAL CAN | SAM |
| A:8260C | 240-59482-B-2-B MS | | 240-212606 | 240-212657 | 12/30/2015 22:09 | 1 | TAL CAN | SAM |
| P:3540C | 240-59482-C-2-C MS | | 240-212723 | 240-212548 | 12/30/2015 07:05 | 1 | TAL CAN | YDN |
| A:8270D | 240-59482-C-2-C MS | | 240-212723 | 240-212548 | 12/31/2015 12:58 | 1 | TAL CAN | MRU |
| P:5030A | 240-59482-C-2-G MS | | 240-212533 | 240-212559 | 12/30/2015 07:39 | 1 | TAL CAN | RTR |
| A:8015C | 240-59482-C-2-G MS | | 240-212533 | 240-212559 | 12/30/2015 10:57 | 1 | TAL CAN | RTR |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Laboratory Chronicle

**Lab ID: 240-59482-2 MSD** **Client ID: F-VB-UST-P-7**

Sample Date/Time: 12/28/2015 10:00 Received Date/Time: 12/29/2015 09:30

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-59482-B-2-C MSD | | 240-212606 | 240-212657 | 12/30/2015 14:25 | 1 | TAL CAN | SAM |
| A:8260C | 240-59482-B-2-C MSD | | 240-212606 | 240-212657 | 12/30/2015 22:31 | 1 | TAL CAN | SAM |
| P:3540C | 240-59482-C-2-D MSD | | 240-212723 | 240-212548 | 12/30/2015 07:05 | 1 | TAL CAN | YDN |
| A:8270D | 240-59482-C-2-D MSD | | 240-212723 | 240-212548 | 12/31/2015 13:23 | 1 | TAL CAN | MRU |
| P:5030A | 240-59482-C-2-H MSD | | 240-212533 | 240-212559 | 12/30/2015 07:39 | 1 | TAL CAN | RTR |
| A:8015C | 240-59482-C-2-H MSD | | 240-212533 | 240-212559 | 12/30/2015 11:31 | 1 | TAL CAN | RTR |

**Lab ID: 240-59482-3** **Client ID: F-VB-UST-P-8**

Sample Date/Time: 12/28/2015 10:30 Received Date/Time: 12/29/2015 09:30

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-59482-B-3-A | | 240-212606 | 240-212657 | 12/30/2015 14:25 | 1 | TAL CAN | SAM |
| A:8260C | 240-59482-B-3-A | | 240-212606 | 240-212657 | 12/30/2015 23:14 | 1 | TAL CAN | SAM |
| P:3540C | 240-59482-C-3-B | | 240-212723 | 240-212548 | 12/30/2015 07:05 | 1 | TAL CAN | YDN |
| A:8270D | 240-59482-C-3-B | | 240-212723 | 240-212548 | 12/31/2015 12:08 | 1 | TAL CAN | MRU |
| P:5030A | 240-59482-C-3-D | | 240-212533 | 240-212559 | 12/30/2015 07:39 | 1 | TAL CAN | RTR |
| A:8015C | 240-59482-C-3-D | | 240-212533 | 240-212559 | 12/30/2015 12:04 | 1 | TAL CAN | RTR |
| P:3540C | 240-59482-C-3-C | | 240-212780 | 240-212549 | 12/30/2015 07:14 | 1 | TAL CAN | YDN |
| A:8015C | 240-59482-C-3-C | | 240-212780 | 240-212549 | 12/31/2015 18:15 | 1 | TAL CAN | HMB |
| P:3540C | 240-59482-C-3-A | | 240-212845 | 240-212544 | 12/30/2015 06:54 | 1 | TAL CAN | YDN |
| A:8082A | 240-59482-C-3-A | | 240-212845 | 240-212544 | 01/04/2016 10:37 | 1 | TAL CAN | HMB |
| A:Moisture | 240-59482-A-3 | | 240-212487 | | 12/29/2015 13:50 | 1 | TAL CAN | BLW |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Laboratory Chronicle

**Lab ID: MB** **Client ID: N/A**

Sample Date/Time: N/A Received Date/Time: N/A

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| A:8260C | MB 240-212606/24 | | 240-212606 | | 12/30/2015 20:00 | 1 | TAL CAN | SAM |
| P:5030C | MB 240-212893/8 | | 240-212893 | | 01/04/2016 11:26 | 1 | TAL CAN | LEE |
| A:8260C | MB 240-212893/8 | | 240-212893 | | 01/04/2016 11:26 | 1 | TAL CAN | LEE |
| P:3540C | MB 240-212548/7-A | | 240-212723 | 240-212548 | 12/30/2015 07:05 | 1 | TAL CAN | YDN |
| A:8270D | MB 240-212548/7-A | | 240-212723 | 240-212548 | 12/31/2015 10:02 | 1 | TAL CAN | MRU |
| P:5030A | MB 240-212559/1-A | | 240-212533 | 240-212559 | 12/30/2015 07:39 | 1 | TAL CAN | RTR |
| A:8015C | MB 240-212559/1-A | | 240-212533 | 240-212559 | 12/30/2015 09:18 | 1 | TAL CAN | RTR |
| P:3540C | MB 240-212549/8-A | | 240-212780 | 240-212549 | 12/30/2015 07:14 | 1 | TAL CAN | YDN |
| A:8015C | MB 240-212549/8-A | | 240-212780 | 240-212549 | 12/31/2015 20:19 | 1 | TAL CAN | HMB |
| P:3540C | MB 240-212544/7-A | | 240-212845 | 240-212544 | 12/30/2015 06:54 | 1 | TAL CAN | YDN |
| A:8082A | MB 240-212544/7-A | | 240-212845 | 240-212544 | 01/04/2016 12:43 | 1 | TAL CAN | HMB |

**Lab ID: LCS** **Client ID: N/A**

Sample Date/Time: N/A Received Date/Time: N/A

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| A:8260C | LCS 240-212606/23 | | 240-212606 | | 12/30/2015 19:39 | 1 | TAL CAN | SAM |
| A:8260C | LCS 240-212606/25 | | 240-212606 | | 12/30/2015 20:22 | 1 | TAL CAN | SAM |
| P:5030C | LCS 240-212893/5 | | 240-212893 | | 01/04/2016 10:18 | 1 | TAL CAN | LEE |
| A:8260C | LCS 240-212893/5 | | 240-212893 | | 01/04/2016 10:18 | 1 | TAL CAN | LEE |
| P:3540C | LCS 240-212548/8-A | | 240-212723 | 240-212548 | 12/30/2015 07:05 | 1 | TAL CAN | YDN |
| A:8270D | LCS 240-212548/8-A | | 240-212723 | 240-212548 | 12/31/2015 10:53 | 1 | TAL CAN | MRU |
| P:5030A | LCS 240-212559/2-A | | 240-212533 | 240-212559 | 12/30/2015 07:39 | 1 | TAL CAN | RTR |
| A:8015C | LCS 240-212559/2-A | | 240-212533 | 240-212559 | 12/30/2015 09:50 | 1 | TAL CAN | RTR |
| P:3540C | LCS 240-212549/9-A | | 240-212780 | 240-212549 | 12/30/2015 07:14 | 1 | TAL CAN | YDN |
| A:8015C | LCS 240-212549/9-A | | 240-212780 | 240-212549 | 12/31/2015 20:50 | 1 | TAL CAN | HMB |
| P:3540C | LCS 240-212544/8-A | | 240-212845 | 240-212544 | 12/30/2015 06:54 | 1 | TAL CAN | YDN |
| A:8082A | LCS 240-212544/8-A | | 240-212845 | 240-212544 | 01/04/2016 13:01 | 1 | TAL CAN | HMB |

**Lab ID: MS** **Client ID: N/A**

Sample Date/Time: 12/29/2015 16:48 Received Date/Time: 12/31/2015 16:25

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-59545-B-2 MS | | 240-212893 | | 01/04/2016 17:00 | 500 | TAL CAN | LEE |
| A:8260C | 240-59545-B-2 MS | | 240-212893 | | 01/04/2016 17:00 | 500 | TAL CAN | LEE |
| P:3540C | 240-59503-G-4-B MS | | 240-212845 | 240-212544 | 12/30/2015 06:54 | 1 | TAL CAN | YDN |
| A:8082A | 240-59503-G-4-B MS | | 240-212845 | 240-212544 | 01/04/2016 12:07 | 1 | TAL CAN | HMB |

Client: Tetra Tech, Inc. Job Number: 240-59482-1

## Laboratory Chronicle

**Lab ID: MSD** **Client ID: N/A**

Sample Date/Time: 12/29/2015 16:48 Received Date/Time: 12/31/2015 16:25

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| P:5030C | 240-59545-B-2 MSD | | 240-212893 | | 01/04/2016 17:23 | 500 | TAL CAN | LEE |
| A:8260C | 240-59545-B-2 MSD | | 240-212893 | | 01/04/2016 17:23 | 500 | TAL CAN | LEE |
| P:3540C | 240-59503-G-4-C MSD | | 240-212845 | 240-212544 | 12/30/2015 06:54 | 1 | TAL CAN | YDN |
| A:8082A | 240-59503-G-4-C MSD | | 240-212845 | 240-212544 | 01/04/2016 12:25 | 1 | TAL CAN | HMB |

**Lab ID: DU** **Client ID: N/A**

Sample Date/Time: 12/16/2015 09:50 Received Date/Time: 12/17/2015 14:07

| Method | Bottle ID | Run | Analysis Batch | Prep Batch | Date Prepared / Analyzed | Dil | Lab | Analyst |
|---|---|---|---|---|---|---|---|---|
| A:Moisture | 240-59255-B-13 DU | | 240-212487 | | 12/29/2015 13:50 | 1 | TAL CAN | BLW |

**Lab References:**
TAL CAN = TestAmerica Canton

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **ex10PPMSPK_00019** | 07/13/17 | 11/23/15 | MEOH, Lot 0000118655 | 1000 mL | ex1016/1260st_00003 | 10 mL | Aroclor-1016 | 10 ug/mL |
| | | | | | | | Aroclor-1260 | 10 ug/mL |
| .ex1016/1260st_00003 | 06/30/21 | | Restek, Lot A0109664 | | (Purchased Reagent) | | Aroclor-1016 | 1000 ug/mL |
| | | | | | | | Aroclor-1260 | 1000 ug/mL |
| **ex2/.2SURRW_00063** | 07/13/17 | 11/23/15 | MEOH, Lot 0000118655 | 2000 mL | EXPESTSURR_00001 | 2 mL | DCB Decachlorobiphenyl | 0.2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.2 ug/mL |
| .EXPESTSURR_00001 | 05/31/21 | | Restek, Lot A0108779 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **exBENZALDEHYD_00033** | 08/31/16 | 12/11/15 | MEOH, Lot 0000118655 | 200 mL | EXLIST1_S11_00001 | 4 mL | Atrazine | 40 ug/mL |
| | | | | | | | Benzaldehyde | 40 ug/mL |
| | | | | | | | Caprolactam | 40 ug/mL |
| .EXLIST1_S11_00001 | 08/31/16 | | Restek, Lot A0108989 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| **exBNASPIKE_00055** | 08/31/16 | 12/07/15 | MEOH, Lot 0000118655 | 200 mL | exLIST1_S1_00004 | 4 mL | 1,1'-Biphenyl | 20 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 20 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 20 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dioxane | 20 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 20 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 40 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 20 ug/mL |
| | | | | | | | 2-Chlorophenol | 20 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2-Methylphenol | 20 ug/mL |
| | | | | | | | 2-Nitroaniline | 20 ug/mL |
| | | | | | | | 2-Nitrophenol | 20 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 20 ug/mL |
| | | | | | | | 3-Methylphenol | 10 ug/mL |
| | | | | | | | 3-Nitroaniline | 20 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 40 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 20 ug/mL |
| | | | | | | | 4-Chloroaniline | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Chlorophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Methylphenol | 10 ug/mL |
| | | | | | | | 4-Nitroaniline | 20 ug/mL |
| | | | | | | | 4-Nitrophenol | 40 ug/mL |
| | | | | | | | Acenaphthene | 20 ug/mL |
| | | | | | | | Acenaphthylene | 20 ug/mL |
| | | | | | | | Acetophenone | 20 ug/mL |
| | | | | | | | Aniline | 20 ug/mL |
| | | | | | | | Anthracene | 20 ug/mL |
| | | | | | | | Azobenzene | 20 ug/mL |
| | | | | | | | Benzo[a]anthracene | 20 ug/mL |
| | | | | | | | Benzo[a]pyrene | 20 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 20 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 20 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 20 ug/mL |
| | | | | | | | Benzyl alcohol | 20 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 20 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 20 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 20 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 20 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 20 ug/mL |
| | | | | | | | Carbazole | 20 ug/mL |
| | | | | | | | Chrysene | 20 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 20 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 20 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 20 ug/mL |
| | | | | | | | Dibenzofuran | 20 ug/mL |
| | | | | | | | Diethyl phthalate | 20 ug/mL |
| | | | | | | | Dimethyl phthalate | 20 ug/mL |
| | | | | | | | Fluoranthene | 20 ug/mL |
| | | | | | | | Fluorene | 20 ug/mL |
| | | | | | | | Hexachlorobenzene | 20 ug/mL |
| | | | | | | | Hexachlorobutadiene | 20 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 20 ug/mL |
| | | | | | | | Hexachloroethane | 20 ug/mL |
| | | | | | | | Hexadecane | 20 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 20 ug/mL |
| | | | | | | | Isophorone | 20 ug/mL |
| | | | | | | | n-Decane | 20 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 40 ug/mL |
| | | | | | | | n-Octadecane | 20 ug/mL |
| | | | | | | | Naphthalene | 20 ug/mL |
| | | | | | | | Nitrobenzene | 20 ug/mL |
| | | | | | | | Pentachlorophenol | 40 ug/mL |
| | | | | | | | Phenanthrene | 20 ug/mL |
| | | | | | | | Phenol | 20 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Pyrene | 20 ug/mL |
| | | | | | | | Pyridine | 20 ug/mL |
| | | | | | EXLIST1_S10_00001 | 4 mL | Benzoic acid | 40 ug/mL |
| | | | | | | | Indene | 40 ug/mL |
| | | | | | exLIST1_S9_00003 | 4 mL | 3,3'-Dichlorobenzidine | 40 ug/mL |
| | | | | | | | Benzidine | 40 ug/mL |
| .exLIST1_S1_00004 | 12/31/16 | | Restek, Lot A0111934 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Methylphenol | 500 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Methylphenol | 500 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| .EXLIST1_S10_00001 | 08/31/16 | | Restek, Lot A0108988 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| .exLIST1_S9_00003 | 01/31/17 | | Restek, Lot A0112567 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| **exBNASURR W_00030** | 07/13/17 | 12/11/15 | MEOH, Lot 0000118655 | 2000 mL | SMLIST1 SURR_00003 | 8 mL | 2,4,6-Tribromophenol (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Nitrobenzene-d5 (Surr) | 20 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 20 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 20 ug/mL |
| .SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **exDIESELSPKw_00069** | 07/06/18 | 11/18/15 | ACETONE, Lot 0000116371 | 200 mL | exDIESELstd_00003 | 10 mL | Diesel | 2521.45 ug/mL |
| .exDIESELstd_00003 | 04/30/22 | | Restek, Lot A0109732 | | (Purchased Reagent) | | Diesel | 50429 ug/mL |
| **exOTPSURRW_00007** | 11/15/16 | 11/23/15 | MEOH, Lot 0000099494 | 1000 mL | exOTPstd_00002 | 20 mL | o-Terphenyl (Surr) | 200 ug/mL |
| .exOTPstd_00002 | 12/31/18 | | Restek, Lot A0110923 | | (Purchased Reagent) | | o-Terphenyl (Surr) | 10000 ug/mL |
| **Gas Surr Cali_00015** | 01/12/16 | 10/12/15 | MeoH, Lot 1823933 | 10 mL | VGGASURROGATE_00032 | 1 mL | Trifluorotoluene (Surr) | 5 ug/mL |
| .VGGASURROGATE_00032 | 08/22/15 | 05/22/15 | MEOH, Lot 49909 | 50 mL | VGSURRSTKGAS_00019 | 1.25 mL | Trifluorotoluene (Surr) | 50 ug/mL |
| ..VGSURRSTKGAS_00019 | 08/31/16 | | ULTRA SCIENTIFIC, Lot CE-1903A | | (Purchased Reagent) | | Trifluorotoluene (Surr) | 2000 ug/mL |
| **SG1248@.05ppm_00018** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 0.5 mL | PCB-1248 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.05 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.1PPM_00015** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 1 mL | PCB-1248 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.1 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.2ppm_00015** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 2 mL | PCB-1248 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.2 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.5ppm_00020** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 5 mL | PCB-1248 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 0.5 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@0.5ppm_00020** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 5 mL | Aroclor-1248 | 0.5 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | Aroclor-1248 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | Aroclor-1248 | 1000 ug/mL |
| **SG1248@1.0ppm_00014** | 04/22/16 | 10/08/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 10 mL | PCB-1248 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1248 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1248@2.0ppm_00014** | 04/22/16 | 08/24/15 | HEXANE, Lot 154006 | 100 mL | SG1248@10ppm_00008 | 20 mL | PCB-1248 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG1248@10ppm_00008 | 08/24/16 | 08/24/15 | Hexane, Lot 151838 | 100 mL | SG1248_00005 | 1 mL | PCB-1248 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 10 ug/mL |
| ..SG1248_00005 | 02/28/21 | | Restek, Lot A0107141 | | (Purchased Reagent) | | PCB-1248 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1248 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1660@.05PPM_00016** | 04/22/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260_00019 | 0.5 mL | PCB-1016 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.05 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.125 mL | DCB Decachlorobiphenyl | 0.0025 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.0025 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | Restek, Lot a109056 | | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | Restek, Lot A0108735 | | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | Restek, Lot A0103185 | | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@0.2ppm_00018** | 04/22/16 | 11/13/15 | HEXANE, Lot 00000101243 | 100 mL | SG1016/1260_00019 | 2 mL | PCB-1016 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.2 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.2 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.5 mL | DCB Decachlorobiphenyl | 0.01 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.01 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@0.5PPM_00058** | 04/22/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260_00019 | 5 mL | PCB-1016 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.5 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 1.25 mL | DCB Decachlorobiphenyl | 0.025 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.025 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@0.5PPM_00058** | 04/22/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260_00019 | 5 mL | Aroclor-1016 | 0.5 ug/mL |
| | | | | | | | Aroclor-1260 | 0.5 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | Aroclor-1016 | 10 ug/mL |
| | | | | | | | Aroclor-1260 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | Aroclor-1016 | 1000 ug/mL |
| | | | | | | | Aroclor-1260 | 1000 ug/mL |
| **SG1660@1.0PPM_00016** | 04/22/16 | 11/13/15 | HEXANE, Lot 00000101243 | 100 mL | SG1016/1260_00019 | 10 mL | PCB-1016 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 2.5 mL | DCB Decachlorobiphenyl | 0.05 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.05 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660@2.0PPM_00015** | 04/22/16 | 11/13/15 | HEXANE, Lot 00000101243 | 100 mL | SG1016/1260_00019 | 20 mL | PCB-1016 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 2 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 5 mL | DCB Decachlorobiphenyl | 0.1 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.1 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **SG1660ICV@1.0_00015** | 04/21/16 | 11/03/15 | HEXANE, Lot 0000101243 | 10 mL | SGPCBIS STOCk_00003 | 0.05 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG1660ICV@1.0_00015** | 04/21/16 | 11/03/15 | HEXANE, Lot 0000101243 | 10 mL | SG1660ICV@100_00005 | 0.1 mL | Aroclor-1016 | 1 ug/mL |
| | | | | | | | Aroclor-1260 | 1 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.25 mL | DCB Decachlorobiphenyl | 0.05 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.05 ug/mL |
| .SG1660ICV@100_00005 | 10/28/16 | 10/28/15 | HEXANE, Lot 0000101243 | 10 mL | SG 1660 ICV_00012 | 1 mL | Aroclor-1016 | 100 ug/mL |
| | | | | | | | Aroclor-1260 | 100 ug/mL |
| ..SG 1660 ICV_00012 | 01/30/22 | | restek, Lot a0114674 | | (Purchased Reagent) | | Aroclor-1016 | 1000 ug/mL |
| | | | | | | | Aroclor-1260 | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG1660STD@0.1_00016** | 04/22/16 | 11/13/15 | HEXANE, Lot 154006 | 100 mL | SG1016/1260_00019 | 1 mL | PCB-1016 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 0.1 ug/mL |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| | | | | | SGPCGSURRSTK_00006 | 0.25 mL | DCB Decachlorobiphenyl | 0.005 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 0.005 ug/mL |
| .SG1016/1260_00019 | 11/13/16 | 11/13/15 | HEXANE, Lot 0000101243 | 100 mL | SG1016/1260MX_00009 | 1 mL | PCB-1016 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1260 Peak 5 | 10 ug/mL |
| ..SG1016/1260MX_00009 | 05/31/21 | | Restek, Lot a109056 | | (Purchased Reagent) | | PCB-1016 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1016 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1260 Peak 4 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1260 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| .SGPCGSURRSTK_00006 | 05/04/16 | 05/04/15 | HEXANE, Lot 0000086800 | 100 mL | SGPESTSURR_00007 | 1 mL | DCB Decachlorobiphenyl | 2 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 2 ug/mL |
| ..SGPESTSURR_00007 | 08/31/20 | | Restek, Lot A0103185 | | (Purchased Reagent) | | DCB Decachlorobiphenyl | 200 ug/mL |
| | | | | | | | Tetrachloro-m-xylene | 200 ug/mL |
| **SG2154@0.05PP_00010** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 0.5 mL | PCB-1221 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.05 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.05 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@0.2PPM_00012** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 2 mL | PCB-1221 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.2 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.2 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@0.5PPM_00021** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 5 mL | PCB-1221 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.5 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.5 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@0.5PPM_00021** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 5 mL | Aroclor-1221 | 0.5 ug/mL |
| | | | | | | | Aroclor-1254 | 0.5 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | Aroclor-1221 | 10 ug/mL |
| | | | | | | | Aroclor-1254 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | Aroclor-1221 | 1000 ug/mL |
| | | | | | | | Aroclor-1254 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **SG2154@1.0PPM_00014** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 10 mL | PCB-1221 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@2.0PPM_00011** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 20 mL | PCB-1221 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 2 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 2 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG2154@O.1PPM_00013** | 03/11/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG2154@10ppm_00009 | 1 mL | PCB-1221 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 0.1 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 0.1 ug/mL |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 0.05 ug/mL |
| .SG2154@10ppm_00009 | 03/11/16 | 06/09/15 | HEXANE, Lot 0000096800 | 100 mL | SG1221/1254_00001 | 1 mL | PCB-1221 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 10 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 10 ug/mL |
| ..SG1221/1254_00001 | 03/11/16 | | restek, Lot a0108261 | | (Purchased Reagent) | | PCB-1221 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1221 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1254 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@.05PPM_00004** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00003 | 0.5 mL | PCB-1232 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 50 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00003 | 09/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/62 STK_00001 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SG32/62 STK_00001 | 04/30/21 | | Restelk, Lot A0108264 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@.2PPM_00005** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 2 mL | PCB-1232 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 200 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@0.1PPM_00006** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 1 mL | PCB-1232 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 100 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@0.5PPM_00010** | 04/22/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00003 | 5 mL | PCB-1232 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 500 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1262 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 500 ug/L |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00003 | 09/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/62 STK_00001 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SG32/62 STK_00001 | 04/30/21 | | Restelk, Lot A0108264 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@0.5PPM_00010** | 04/22/16 | 10/06/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00003 | 5 mL | Aroclor-1232 | 500 ug/L |
| | | | | | | | PCB-1262 | 500 ug/L |
| .SG32/6210PPM_00003 | 09/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/62 STK_00001 | 1 mL | Aroclor-1232 | 10000 ug/L |
| | | | | | | | PCB-1262 | 10000 ug/L |
| ..SG32/62 STK_00001 | 04/30/21 | | Restelk, Lot A0108264 | | (Purchased Reagent) | | Aroclor-1232 | 1000 ug/mL |
| | | | | | | | PCB-1262 | 1000 ug/mL |
| **SG3262@1.0PPM_00005** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 10 mL | PCB-1232 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG3262@2.0ppm_00004** | 03/25/16 | 09/25/15 | HEXANE, Lot 154006 | 100 mL | SG32/6210PPM_00002 | 20 mL | PCB-1232 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 2000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG32/6210PPM_00002 | 04/24/16 | 04/24/15 | HEXANE, Lot 0000096800 | 100 mL | SGPCB1232STD_00005 | 1 mL | PCB-1232 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1232 Peak 5 | 10000 ug/L |
| | | | | | SGPCB1262STD_00005 | 1 mL | PCB-1262 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1262 Peak 5 | 10000 ug/L |
| ..SGPCB1232STD_00005 | 04/30/18 | | restek, Lot A086119 | | (Purchased Reagent) | | PCB-1232 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1232 Peak 5 | 1000 ug/mL |
| ..SGPCB1262STD_00005 | 05/31/20 | | restek, Lot A0101379 | | (Purchased Reagent) | | PCB-1262 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1262 Peak 4 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1262 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG42/68@1.0PP_00008** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 10 mL | PCB-1242 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@.05PPM_00005** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 0.5 mL | PCB-1242 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 50 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 50 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@.1PPM_00006** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 1 mL | PCB-1242 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 100 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 100 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@.2PPM_00005** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 2 mL | PCB-1242 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 200 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 200 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@0.5PPM_00007** | 12/24/15 | 06/24/15 | HEXANE, Lot 0000086560 | 100 mL | SG42/6810PPM_00002 | 5 mL | PCB-1242 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 500 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1268 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 500 ug/L |
| | | | | | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@0.5PPM_00011** | 04/07/16 | 10/07/15 | HEXANE, Lot 154006 | 100 mL | SGPCBIS STOCk_00002 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SGPCBIS STOCk_00002 | 04/22/16 | 04/22/15 | HEXANE, Lot 0000086560 | 100 mL | ISTD_00002 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00002 | 04/30/18 | | Restek, Lot A0110081 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SG4268@0.5PPM_00011** | 04/07/16 | 10/07/15 | HEXANE, Lot 154006 | 100 mL | SG42/6810PPM_00003 | 5 mL | Aroclor-1242 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 500 ug/L |
| | | | | | | | PCB-1268 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 500 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 500 ug/L |
| .SG42/6810PPM_00003 | 10/07/16 | 10/07/15 | HEXANE, Lot 154006 | 100 mL | SG 42/68 STK_00001 | 1 mL | Aroclor-1242 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | | | PCB-1268 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG 42/68 STK_00001 | 04/30/21 | | RESTEK, Lot A0108269 | | (Purchased Reagent) | | Aroclor-1242 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| | | | | | | | PCB-1268 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| **SG4268@2.0PPM_00004** | 04/01/16 | 10/01/15 | HEXANE, Lot 00000154006 | 100 mL | SG42/6810PPM_00002 | 20 mL | PCB-1242 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 1 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 2000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 2000 ug/L |
| | | | | | SGPCBIS STOCk_00003 | 0.5 mL | 1-Bromo-2-nitrobenzene | 50 ug/L |
| .SG42/6810PPM_00002 | 04/28/16 | 04/28/15 | HEXANE, Lot 0000096800 | 100 mL | SG1242_00005 | 1 mL | PCB-1242 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1242 Peak 5 | 10000 ug/L |
| | | | | | SGpcb1268std_00006 | 1 mL | PCB-1268 Peak 1 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 2 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 3 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 4 | 10000 ug/L |
| | | | | | | | PCB-1268 Peak 5 | 10000 ug/L |
| ..SG1242_00005 | 11/30/18 | | Restek, Lot A090182 | | (Purchased Reagent) | | PCB-1242 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 2 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1242 Peak 5 | 1000 ug/mL |
| ..SGpcb1268std_00006 | 12/31/19 | | Restek, Lot A097795 | | (Purchased Reagent) | | PCB-1268 Peak 1 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 2 | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | PCB-1268 Peak 3 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 4 | 1000 ug/mL |
| | | | | | | | PCB-1268 Peak 5 | 1000 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SGDROCAL_00021** | 05/30/16 | 11/30/15 | MECL2, Lot SGDRODCM_00001 | 25 mL | SGDROCALSTK_00020 | 5 mL | Diesel | 10000 ug/mL |
| | | | | | SGDROSURROTP_00027 | 2 mL | o-Terphenyl (Surr) | 800 ug/mL |
| .SGDROCALSTK_00020 | 04/30/22 | | Restek, Lot A0109732 | | (Purchased Reagent) | | Diesel | 50000 ug/mL |
| .SGDROSURROTP_00027 | 12/31/18 | | Restek, Lot A0110923 | | (Purchased Reagent) | | o-Terphenyl (Surr) | 10000 ug/mL |
| **SGDROICV_00015** | 08/26/16 | 08/26/15 | MECL2, Lot 2091477 | 25 mL | SGDROICVSTK_00008 | 1 mL | Diesel | 2006 ug/mL |
| .SGDROICVSTK_00008 | 08/31/18 | | Ultra, Lot CL-2929 | | (Purchased Reagent) | | Diesel | 50150 ug/mL |
| **SGPCBISTD_00003** | 04/22/16 | 10/13/15 | HEXANE, Lot 154006 | 100 mL | SGPCBIS STOCk_00003 | 10 mL | 1-Bromo-2-nitrobenzene | 1 ug/mL |
| .SGPCBIS STOCk_00003 | 04/22/16 | 04/22/15 | HEXANE, Lot 000017549 | 100 mL | ISTD_00001 | 1 mL | 1-Bromo-2-nitrobenzene | 10 ug/mL |
| ..ISTD_00001 | 04/30/18 | | Restek, Lot A0108735 | | (Purchased Reagent) | | 1-Bromo-2-nitrobenzene | 1000 ug/mL |
| **SMIS80PPMW_00013** | 07/10/16 | 07/10/15 | MECL2, Lot 0000103549 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| .SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| **SMLIST1 L1 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1PAH ST_00004 | 2 uL | 1,1'-Biphenyl | 0.1 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 0.1 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 0.1 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 0.1 ug/mL |
| | | | | | | | Acenaphthene | 0.1 ug/mL |
| | | | | | | | Acenaphthylene | 0.1 ug/mL |
| | | | | | | | Anthracene | 0.1 ug/mL |
| | | | | | | | Benzo[a]anthracene | 0.1 ug/mL |
| | | | | | | | Benzo[a]pyrene | 0.1 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 0.1 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 0.1 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 0.1 ug/mL |
| | | | | | | | Chrysene | 0.1 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Dibenz(a,h)anthracene | 0.1 ug/mL |
| | | | | | | | Dibenzofuran | 0.1 ug/mL |
| | | | | | | | Fluoranthene | 0.1 ug/mL |
| | | | | | | | Fluorene | 0.1 ug/mL |
| | | | | | | | Hexachlorobenzene | 0.1 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 0.1 ug/mL |
| | | | | | | | Naphthalene | 0.1 ug/mL |
| | | | | | | | Phenanthrene | 0.1 ug/mL |
| | | | | | | | Pyrene | 0.1 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 0.1 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 0.1 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 0.1 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1PAH ST_00004 | 05/31/16 | 01/26/15 | MECL2, Lot 0000093871 | 3 mL | SMLIST1 PAH_00004 | 300 uL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 60 uL | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 PAH_00004 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L2 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 10 uL | 1,1'-Biphenyl | 0.5 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 0.5 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 0.5 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 0.5 ug/mL |
| | | | | | | | 1,4-Dioxane | 0.5 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 0.5 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 0.5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,4-Dinitrophenol | 1 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 0.5 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 0.5 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 0.5 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 0.5 ug/mL |
| | | | | | | | 2-Chlorophenol | 0.5 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 0.5 ug/mL |
| | | | | | | | 2-Methylphenol | 0.5 ug/mL |
| | | | | | | | 2-Nitroaniline | 0.5 ug/mL |
| | | | | | | | 2-Nitrophenol | 0.5 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 0.5 ug/mL |
| | | | | | | | 3-Nitroaniline | 0.5 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 1 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 0.5 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 0.5 ug/mL |
| | | | | | | | 4-Chloroaniline | 0.5 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 0.5 ug/mL |
| | | | | | | | 4-Nitroaniline | 0.5 ug/mL |
| | | | | | | | 4-Nitrophenol | 1 ug/mL |
| | | | | | | | Acenaphthene | 0.5 ug/mL |
| | | | | | | | Acenaphthylene | 0.5 ug/mL |
| | | | | | | | Acetophenone | 0.5 ug/mL |
| | | | | | | | Aniline | 0.5 ug/mL |
| | | | | | | | Anthracene | 0.5 ug/mL |
| | | | | | | | Azobenzene | 0.5 ug/mL |
| | | | | | | | Benzo[a]anthracene | 0.5 ug/mL |
| | | | | | | | Benzo[a]pyrene | 0.5 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 0.5 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 0.5 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 0.5 ug/mL |
| | | | | | | | Benzyl alcohol | 0.5 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 0.5 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 0.5 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 0.5 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 0.5 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 0.5 ug/mL |
| | | | | | | | Carbazole | 0.5 ug/mL |
| | | | | | | | Chrysene | 0.5 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 0.5 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 0.5 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 0.5 ug/mL |
| | | | | | | | Dibenzofuran | 0.5 ug/mL |
| | | | | | | | Diethyl phthalate | 0.5 ug/mL |
| | | | | | | | Dimethyl phthalate | 0.5 ug/mL |
| | | | | | | | Fluoranthene | 0.5 ug/mL |
| | | | | | | | Fluorene | 0.5 ug/mL |
| | | | | | | | Hexachlorobenzene | 0.5 ug/mL |
| | | | | | | | Hexachlorobutadiene | 0.5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Hexachlorocyclopentadiene | 0.5 ug/mL |
| | | | | | | | Hexachloroethane | 0.5 ug/mL |
| | | | | | | | Hexadecane | 0.5 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 0.5 ug/mL |
| | | | | | | | Isophorone | 0.5 ug/mL |
| | | | | | | | n-Decane | 0.5 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 0.5 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 0.5 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 1 ug/mL |
| | | | | | | | n-Octadecane | 0.5 ug/mL |
| | | | | | | | Naphthalene | 0.5 ug/mL |
| | | | | | | | Nitrobenzene | 0.5 ug/mL |
| | | | | | | | Pentachlorophenol | 1 ug/mL |
| | | | | | | | Phenanthrene | 0.5 ug/mL |
| | | | | | | | Phenol | 0.5 ug/mL |
| | | | | | | | Pyrene | 0.5 ug/mL |
| | | | | | | | Pyridine | 0.5 ug/mL |
| | | | | | | | Benzoic acid | 1 ug/mL |
| | | | | | | | Indene | 1 ug/mL |
| | | | | | | | Atrazine | 1 ug/mL |
| | | | | | | | Benzaldehyde | 1 ug/mL |
| | | | | | | | Caprolactam | 1 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 1 ug/mL |
| | | | | | | | Benzidine | 1 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 0.5 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 0.5 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 0.5 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 0.5 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 0.5 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 0.5 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L3 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 20 uL | 1,1'-Biphenyl | 1 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,3-Dichlorobenzene | 1 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1 ug/mL |
| | | | | | | | 1,4-Dioxane | 1 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1 ug/mL |
| | | | | | | | 2-Chlorophenol | 1 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1 ug/mL |
| | | | | | | | 2-Methylphenol | 1 ug/mL |
| | | | | | | | 2-Nitroaniline | 1 ug/mL |
| | | | | | | | 2-Nitrophenol | 1 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1 ug/mL |
| | | | | | | | 3-Nitroaniline | 1 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1 ug/mL |
| | | | | | | | 4-Chloroaniline | 1 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1 ug/mL |
| | | | | | | | 4-Nitroaniline | 1 ug/mL |
| | | | | | | | 4-Nitrophenol | 2 ug/mL |
| | | | | | | | Acenaphthene | 1 ug/mL |
| | | | | | | | Acenaphthylene | 1 ug/mL |
| | | | | | | | Acetophenone | 1 ug/mL |
| | | | | | | | Aniline | 1 ug/mL |
| | | | | | | | Anthracene | 1 ug/mL |
| | | | | | | | Azobenzene | 1 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1 ug/mL |
| | | | | | | | Benzyl alcohol | 1 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1 ug/mL |
| | | | | | | | Carbazole | 1 ug/mL |
| | | | | | | | Chrysene | 1 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Di-n-butyl phthalate | 1 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1 ug/mL |
| | | | | | | | Dibenzofuran | 1 ug/mL |
| | | | | | | | Diethyl phthalate | 1 ug/mL |
| | | | | | | | Dimethyl phthalate | 1 ug/mL |
| | | | | | | | Fluoranthene | 1 ug/mL |
| | | | | | | | Fluorene | 1 ug/mL |
| | | | | | | | Hexachlorobenzene | 1 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1 ug/mL |
| | | | | | | | Hexachloroethane | 1 ug/mL |
| | | | | | | | Hexadecane | 1 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1 ug/mL |
| | | | | | | | Isophorone | 1 ug/mL |
| | | | | | | | n-Decane | 1 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2 ug/mL |
| | | | | | | | n-Octadecane | 1 ug/mL |
| | | | | | | | Naphthalene | 1 ug/mL |
| | | | | | | | Nitrobenzene | 1 ug/mL |
| | | | | | | | Pentachlorophenol | 2 ug/mL |
| | | | | | | | Phenanthrene | 1 ug/mL |
| | | | | | | | Phenol | 1 ug/mL |
| | | | | | | | Pyrene | 1 ug/mL |
| | | | | | | | Pyridine | 1 ug/mL |
| | | | | | | | Benzoic acid | 2 ug/mL |
| | | | | | | | Indene | 2 ug/mL |
| | | | | | | | Atrazine | 2 ug/mL |
| | | | | | | | Benzaldehyde | 2 ug/mL |
| | | | | | | | Caprolactam | 2 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 2 ug/mL |
| | | | | | | | Benzidine | 2 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 1 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 1 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 1 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 1 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 1 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 1 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | Restek, Lot A0101020 | | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L4 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 40 uL | 1,1'-Biphenyl | 2 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 2 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 2 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 2 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2 ug/mL |
| | | | | | | | 1,4-Dioxane | 2 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 2 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 2 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 2 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 2 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 2 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 2 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 4 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 2 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 2 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 2 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 2 ug/mL |
| | | | | | | | 2-Chlorophenol | 2 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 2 ug/mL |
| | | | | | | | 2-Methylphenol | 2 ug/mL |
| | | | | | | | 2-Nitroaniline | 2 ug/mL |
| | | | | | | | 2-Nitrophenol | 2 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 2 ug/mL |
| | | | | | | | 3-Nitroaniline | 2 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 4 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 2 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 2 ug/mL |
| | | | | | | | 4-Chloroaniline | 2 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 2 ug/mL |
| | | | | | | | 4-Nitroaniline | 2 ug/mL |
| | | | | | | | 4-Nitrophenol | 4 ug/mL |
| | | | | | | | Acenaphthene | 2 ug/mL |
| | | | | | | | Acenaphthylene | 2 ug/mL |
| | | | | | | | Acetophenone | 2 ug/mL |
| | | | | | | | Aniline | 2 ug/mL |
| | | | | | | | Anthracene | 2 ug/mL |
| | | | | | | | Azobenzene | 2 ug/mL |
| | | | | | | | Benzo[a]anthracene | 2 ug/mL |
| | | | | | | | Benzo[a]pyrene | 2 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 2 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[g,h,i]perylene | 2 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 2 ug/mL |
| | | | | | | | Benzyl alcohol | 2 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 2 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 2 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 2 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 2 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 2 ug/mL |
| | | | | | | | Carbazole | 2 ug/mL |
| | | | | | | | Chrysene | 2 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 2 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 2 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 2 ug/mL |
| | | | | | | | Dibenzofuran | 2 ug/mL |
| | | | | | | | Diethyl phthalate | 2 ug/mL |
| | | | | | | | Dimethyl phthalate | 2 ug/mL |
| | | | | | | | Fluoranthene | 2 ug/mL |
| | | | | | | | Fluorene | 2 ug/mL |
| | | | | | | | Hexachlorobenzene | 2 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 2 ug/mL |
| | | | | | | | Hexachloroethane | 2 ug/mL |
| | | | | | | | Hexadecane | 2 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 2 ug/mL |
| | | | | | | | Isophorone | 2 ug/mL |
| | | | | | | | n-Decane | 2 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 2 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 2 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 4 ug/mL |
| | | | | | | | n-Octadecane | 2 ug/mL |
| | | | | | | | Naphthalene | 2 ug/mL |
| | | | | | | | Nitrobenzene | 2 ug/mL |
| | | | | | | | Pentachlorophenol | 4 ug/mL |
| | | | | | | | Phenanthrene | 2 ug/mL |
| | | | | | | | Phenol | 2 ug/mL |
| | | | | | | | Pyrene | 2 ug/mL |
| | | | | | | | Pyridine | 2 ug/mL |
| | | | | | | | Benzoic acid | 4 ug/mL |
| | | | | | | | Indene | 4 ug/mL |
| | | | | | | | Atrazine | 4 ug/mL |
| | | | | | | | Benzaldehyde | 4 ug/mL |
| | | | | | | | Caprolactam | 4 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 4 ug/mL |
| | | | | | | | Benzidine | 4 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 2 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 2 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 2 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 2 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Phenol-d5 (Surr) | 2 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 2 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L5 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 100 uL | 1,1'-Biphenyl | 5 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 5 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 5 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 5 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 5 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 5 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 5 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 5 ug/mL |
| | | | | | | | 1,4-Dioxane | 5 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 5 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 5 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 5 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 5 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 5 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 5 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 10 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 5 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 5 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 5 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 5 ug/mL |
| | | | | | | | 2-Chlorophenol | 5 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 5 ug/mL |
| | | | | | | | 2-Methylphenol | 5 ug/mL |
| | | | | | | | 2-Nitroaniline | 5 ug/mL |
| | | | | | | | 2-Nitrophenol | 5 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 5 ug/mL |
| | | | | | | | 3-Nitroaniline | 5 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 10 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 5 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 5 ug/mL |
| | | | | | | | 4-Chloroaniline | 5 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 5 ug/mL |
| | | | | | | | 4-Nitroaniline | 5 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Nitrophenol | 10 ug/mL |
| | | | | | | | Acenaphthene | 5 ug/mL |
| | | | | | | | Acenaphthylene | 5 ug/mL |
| | | | | | | | Acetophenone | 5 ug/mL |
| | | | | | | | Aniline | 5 ug/mL |
| | | | | | | | Anthracene | 5 ug/mL |
| | | | | | | | Azobenzene | 5 ug/mL |
| | | | | | | | Benzo[a]anthracene | 5 ug/mL |
| | | | | | | | Benzo[a]pyrene | 5 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 5 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 5 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 5 ug/mL |
| | | | | | | | Benzyl alcohol | 5 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 5 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 5 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 5 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 5 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 5 ug/mL |
| | | | | | | | Carbazole | 5 ug/mL |
| | | | | | | | Chrysene | 5 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 5 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 5 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 5 ug/mL |
| | | | | | | | Dibenzofuran | 5 ug/mL |
| | | | | | | | Diethyl phthalate | 5 ug/mL |
| | | | | | | | Dimethyl phthalate | 5 ug/mL |
| | | | | | | | Fluoranthene | 5 ug/mL |
| | | | | | | | Fluorene | 5 ug/mL |
| | | | | | | | Hexachlorobenzene | 5 ug/mL |
| | | | | | | | Hexachlorobutadiene | 5 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 5 ug/mL |
| | | | | | | | Hexachloroethane | 5 ug/mL |
| | | | | | | | Hexadecane | 5 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 5 ug/mL |
| | | | | | | | Isophorone | 5 ug/mL |
| | | | | | | | n-Decane | 5 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 5 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 5 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 10 ug/mL |
| | | | | | | | n-Octadecane | 5 ug/mL |
| | | | | | | | Naphthalene | 5 ug/mL |
| | | | | | | | Nitrobenzene | 5 ug/mL |
| | | | | | | | Pentachlorophenol | 10 ug/mL |
| | | | | | | | Phenanthrene | 5 ug/mL |
| | | | | | | | Phenol | 5 ug/mL |
| | | | | | | | Pyrene | 5 ug/mL |
| | | | | | | | Pyridine | 5 ug/mL |
| | | | | | | | Benzoic acid | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Indene | 10 ug/mL |
| | | | | | | | Atrazine | 10 ug/mL |
| | | | | | | | Benzaldehyde | 10 ug/mL |
| | | | | | | | Caprolactam | 10 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 10 ug/mL |
| | | | | | | | Benzidine | 10 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 5 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L6 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 200 uL | 1,1'-Biphenyl | 10 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 10 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 10 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 10 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 10 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 10 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 10 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 10 ug/mL |
| | | | | | | | 1,4-Dioxane | 10 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 10 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 10 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 10 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 10 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 10 ug/mL |
| | | | | | | | 2-Chlorophenol | 10 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 10 ug/mL |
| | | | | | | | 2-Methylphenol | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Nitroaniline | 10 ug/mL |
| | | | | | | | 2-Nitrophenol | 10 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 10 ug/mL |
| | | | | | | | 3-Nitroaniline | 10 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 20 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 10 ug/mL |
| | | | | | | | 4-Chloroaniline | 10 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Nitroaniline | 10 ug/mL |
| | | | | | | | 4-Nitrophenol | 20 ug/mL |
| | | | | | | | Acenaphthene | 10 ug/mL |
| | | | | | | | Acenaphthylene | 10 ug/mL |
| | | | | | | | Acetophenone | 10 ug/mL |
| | | | | | | | Aniline | 10 ug/mL |
| | | | | | | | Anthracene | 10 ug/mL |
| | | | | | | | Azobenzene | 10 ug/mL |
| | | | | | | | Benzo[a]anthracene | 10 ug/mL |
| | | | | | | | Benzo[a]pyrene | 10 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 10 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 10 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 10 ug/mL |
| | | | | | | | Benzyl alcohol | 10 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 10 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 10 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 10 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 10 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 10 ug/mL |
| | | | | | | | Carbazole | 10 ug/mL |
| | | | | | | | Chrysene | 10 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 10 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 10 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 10 ug/mL |
| | | | | | | | Dibenzofuran | 10 ug/mL |
| | | | | | | | Diethyl phthalate | 10 ug/mL |
| | | | | | | | Dimethyl phthalate | 10 ug/mL |
| | | | | | | | Fluoranthene | 10 ug/mL |
| | | | | | | | Fluorene | 10 ug/mL |
| | | | | | | | Hexachlorobenzene | 10 ug/mL |
| | | | | | | | Hexachlorobutadiene | 10 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 10 ug/mL |
| | | | | | | | Hexachloroethane | 10 ug/mL |
| | | | | | | | Hexadecane | 10 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 10 ug/mL |
| | | | | | | | Isophorone | 10 ug/mL |
| | | | | | | | n-Decane | 10 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 10 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | N-Nitrosodiphenylamine | 20 ug/mL |
| | | | | | | | n-Octadecane | 10 ug/mL |
| | | | | | | | Naphthalene | 10 ug/mL |
| | | | | | | | Nitrobenzene | 10 ug/mL |
| | | | | | | | Pentachlorophenol | 20 ug/mL |
| | | | | | | | Phenanthrene | 10 ug/mL |
| | | | | | | | Phenol | 10 ug/mL |
| | | | | | | | Pyrene | 10 ug/mL |
| | | | | | | | Pyridine | 10 ug/mL |
| | | | | | | | Benzoic acid | 20 ug/mL |
| | | | | | | | Indene | 20 ug/mL |
| | | | | | | | Atrazine | 20 ug/mL |
| | | | | | | | Benzaldehyde | 20 ug/mL |
| | | | | | | | Caprolactam | 20 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 20 ug/mL |
| | | | | | | | Benzidine | 20 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 10 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 10 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 10 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 10 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L7 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 300 uL | 1,1'-Biphenyl | 15 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 15 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 15 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 15 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 15 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 15 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 15 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 15 ug/mL |
| | | | | | | | 1,4-Dioxane | 15 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 15 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 15 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 15 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 15 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2,4-Dichlorophenol | 15 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 15 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 30 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 15 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 15 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 15 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 15 ug/mL |
| | | | | | | | 2-Chlorophenol | 15 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 15 ug/mL |
| | | | | | | | 2-Methylphenol | 15 ug/mL |
| | | | | | | | 2-Nitroaniline | 15 ug/mL |
| | | | | | | | 2-Nitrophenol | 15 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 15 ug/mL |
| | | | | | | | 3-Nitroaniline | 15 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 30 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 15 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 15 ug/mL |
| | | | | | | | 4-Chloroaniline | 15 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 15 ug/mL |
| | | | | | | | 4-Nitroaniline | 15 ug/mL |
| | | | | | | | 4-Nitrophenol | 30 ug/mL |
| | | | | | | | Acenaphthene | 15 ug/mL |
| | | | | | | | Acenaphthylene | 15 ug/mL |
| | | | | | | | Acetophenone | 15 ug/mL |
| | | | | | | | Aniline | 15 ug/mL |
| | | | | | | | Anthracene | 15 ug/mL |
| | | | | | | | Azobenzene | 15 ug/mL |
| | | | | | | | Benzo[a]anthracene | 15 ug/mL |
| | | | | | | | Benzo[a]pyrene | 15 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 15 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 15 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 15 ug/mL |
| | | | | | | | Benzyl alcohol | 15 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 15 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 15 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 15 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 15 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 15 ug/mL |
| | | | | | | | Carbazole | 15 ug/mL |
| | | | | | | | Chrysene | 15 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 15 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 15 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 15 ug/mL |
| | | | | | | | Dibenzofuran | 15 ug/mL |
| | | | | | | | Diethyl phthalate | 15 ug/mL |
| | | | | | | | Dimethyl phthalate | 15 ug/mL |
| | | | | | | | Fluoranthene | 15 ug/mL |
| | | | | | | | Fluorene | 15 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Hexachlorobenzene | 15 ug/mL |
| | | | | | | | Hexachlorobutadiene | 15 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 15 ug/mL |
| | | | | | | | Hexachloroethane | 15 ug/mL |
| | | | | | | | Hexadecane | 15 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 15 ug/mL |
| | | | | | | | Isophorone | 15 ug/mL |
| | | | | | | | n-Decane | 15 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 15 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 15 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 30 ug/mL |
| | | | | | | | n-Octadecane | 15 ug/mL |
| | | | | | | | Naphthalene | 15 ug/mL |
| | | | | | | | Nitrobenzene | 15 ug/mL |
| | | | | | | | Pentachlorophenol | 30 ug/mL |
| | | | | | | | Phenanthrene | 15 ug/mL |
| | | | | | | | Phenol | 15 ug/mL |
| | | | | | | | Pyrene | 15 ug/mL |
| | | | | | | | Pyridine | 15 ug/mL |
| | | | | | | | Benzoic acid | 30 ug/mL |
| | | | | | | | Indene | 30 ug/mL |
| | | | | | | | Atrazine | 30 ug/mL |
| | | | | | | | Benzaldehyde | 30 ug/mL |
| | | | | | | | Caprolactam | 30 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 30 ug/mL |
| | | | | | | | Benzidine | 30 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 15 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 15 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 15 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 15 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 15 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 15 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 L8 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 400 uL | 1,1'-Biphenyl | 20 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 20 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 20 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 20 ug/mL |
| | | | | | | | 1,4-Dioxane | 20 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 20 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 40 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 20 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 20 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 20 ug/mL |
| | | | | | | | 2-Chlorophenol | 20 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 20 ug/mL |
| | | | | | | | 2-Methylphenol | 20 ug/mL |
| | | | | | | | 2-Nitroaniline | 20 ug/mL |
| | | | | | | | 2-Nitrophenol | 20 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 20 ug/mL |
| | | | | | | | 3-Nitroaniline | 20 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 40 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 20 ug/mL |
| | | | | | | | 4-Chloroaniline | 20 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 20 ug/mL |
| | | | | | | | 4-Nitroaniline | 20 ug/mL |
| | | | | | | | 4-Nitrophenol | 40 ug/mL |
| | | | | | | | Acenaphthene | 20 ug/mL |
| | | | | | | | Acenaphthylene | 20 ug/mL |
| | | | | | | | Acetophenone | 20 ug/mL |
| | | | | | | | Aniline | 20 ug/mL |
| | | | | | | | Anthracene | 20 ug/mL |
| | | | | | | | Azobenzene | 20 ug/mL |
| | | | | | | | Benzo[a]anthracene | 20 ug/mL |
| | | | | | | | Benzo[a]pyrene | 20 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 20 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 20 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 20 ug/mL |
| | | | | | | | Benzyl alcohol | 20 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 20 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 20 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 20 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 20 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 20 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Carbazole | 20 ug/mL |
| | | | | | | | Chrysene | 20 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 20 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 20 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 20 ug/mL |
| | | | | | | | Dibenzofuran | 20 ug/mL |
| | | | | | | | Diethyl phthalate | 20 ug/mL |
| | | | | | | | Dimethyl phthalate | 20 ug/mL |
| | | | | | | | Fluoranthene | 20 ug/mL |
| | | | | | | | Fluorene | 20 ug/mL |
| | | | | | | | Hexachlorobenzene | 20 ug/mL |
| | | | | | | | Hexachlorobutadiene | 20 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 20 ug/mL |
| | | | | | | | Hexachloroethane | 20 ug/mL |
| | | | | | | | Hexadecane | 20 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 20 ug/mL |
| | | | | | | | Isophorone | 20 ug/mL |
| | | | | | | | n-Decane | 20 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 20 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 40 ug/mL |
| | | | | | | | n-Octadecane | 20 ug/mL |
| | | | | | | | Naphthalene | 20 ug/mL |
| | | | | | | | Nitrobenzene | 20 ug/mL |
| | | | | | | | Pentachlorophenol | 40 ug/mL |
| | | | | | | | Phenanthrene | 20 ug/mL |
| | | | | | | | Phenol | 20 ug/mL |
| | | | | | | | Pyrene | 20 ug/mL |
| | | | | | | | Pyridine | 20 ug/mL |
| | | | | | | | Benzoic acid | 40 ug/mL |
| | | | | | | | Indene | 40 ug/mL |
| | | | | | | | Atrazine | 40 ug/mL |
| | | | | | | | Benzaldehyde | 40 ug/mL |
| | | | | | | | Caprolactam | 40 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 40 ug/mL |
| | | | | | | | Benzidine | 40 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 20 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 20 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 20 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 20 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 20 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **SMLIST1 L9 W_00006** | 04/20/16 | 04/21/15 | MECL2, Lot 0000103549 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| | | | | | SMLIST1 STOCK_00006 | 500 uL | 1,1'-Biphenyl | 25 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 25 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 25 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 25 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 25 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 25 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 25 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 25 ug/mL |
| | | | | | | | 1,4-Dioxane | 25 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 25 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 25 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 25 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 25 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 25 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 25 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 50 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 25 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 25 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 25 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 25 ug/mL |
| | | | | | | | 2-Chlorophenol | 25 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 25 ug/mL |
| | | | | | | | 2-Methylphenol | 25 ug/mL |
| | | | | | | | 2-Nitroaniline | 25 ug/mL |
| | | | | | | | 2-Nitrophenol | 25 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 25 ug/mL |
| | | | | | | | 3-Nitroaniline | 25 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 50 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 25 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 25 ug/mL |
| | | | | | | | 4-Chloroaniline | 25 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 25 ug/mL |
| | | | | | | | 4-Nitroaniline | 25 ug/mL |
| | | | | | | | 4-Nitrophenol | 50 ug/mL |
| | | | | | | | Acenaphthene | 25 ug/mL |
| | | | | | | | Acenaphthylene | 25 ug/mL |
| | | | | | | | Acetophenone | 25 ug/mL |
| | | | | | | | Aniline | 25 ug/mL |
| | | | | | | | Anthracene | 25 ug/mL |
| | | | | | | | Azobenzene | 25 ug/mL |
| | | | | | | | Benzo[a]anthracene | 25 ug/mL |
| | | | | | | | Benzo[a]pyrene | 25 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Benzo[b]fluoranthene | 25 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 25 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 25 ug/mL |
| | | | | | | | Benzyl alcohol | 25 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 25 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 25 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 25 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 25 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 25 ug/mL |
| | | | | | | | Carbazole | 25 ug/mL |
| | | | | | | | Chrysene | 25 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 25 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 25 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 25 ug/mL |
| | | | | | | | Dibenzofuran | 25 ug/mL |
| | | | | | | | Diethyl phthalate | 25 ug/mL |
| | | | | | | | Dimethyl phthalate | 25 ug/mL |
| | | | | | | | Fluoranthene | 25 ug/mL |
| | | | | | | | Fluorene | 25 ug/mL |
| | | | | | | | Hexachlorobenzene | 25 ug/mL |
| | | | | | | | Hexachlorobutadiene | 25 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 25 ug/mL |
| | | | | | | | Hexachloroethane | 25 ug/mL |
| | | | | | | | Hexadecane | 25 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 25 ug/mL |
| | | | | | | | Isophorone | 25 ug/mL |
| | | | | | | | n-Decane | 25 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 25 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 25 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 50 ug/mL |
| | | | | | | | n-Octadecane | 25 ug/mL |
| | | | | | | | Naphthalene | 25 ug/mL |
| | | | | | | | Nitrobenzene | 25 ug/mL |
| | | | | | | | Pentachlorophenol | 50 ug/mL |
| | | | | | | | Phenanthrene | 25 ug/mL |
| | | | | | | | Phenol | 25 ug/mL |
| | | | | | | | Pyrene | 25 ug/mL |
| | | | | | | | Pyridine | 25 ug/mL |
| | | | | | | | Benzoic acid | 50 ug/mL |
| | | | | | | | Indene | 50 ug/mL |
| | | | | | | | Atrazine | 50 ug/mL |
| | | | | | | | Benzaldehyde | 50 ug/mL |
| | | | | | | | Caprolactam | 50 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 50 ug/mL |
| | | | | | | | Benzidine | 50 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 25 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 25 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 25 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Nitrobenzene-d5 (Surr) | 25 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 25 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 25 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| .SMLIST1 STOCK_00006 | 05/31/16 | 04/20/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 S1_00005 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 100 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 100 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 100 ug/mL |
| | | | | | | | 1,4-Dioxane | 100 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Aniline | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Azobenzene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | Benzyl alcohol | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Hexadecane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | n-Decane | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | n-Octadecane | 100 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | | | Pyridine | 100 ug/mL |
| | | | | | SMLIST1 S10_00001 | 1 mL | Benzoic acid | 200 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Indene | 200 ug/mL |
| | | | | | SMLIST1 S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1 S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | | | Benzidine | 200 ug/mL |
| | | | | | SMLIST1 SURR_00003 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 S1_00005 | 05/31/16 | | Restek, Lot A0107399 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 1,2,4,5-Tetrachlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,2-Diphenylhydrazine | 1000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,3-Dinitrobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 1000 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 1-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2,3,4,6-Tetrachlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Aniline | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Azobenzene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzyl alcohol | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Hexadecane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | n-Decane | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodimethylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | n-Octadecane | 1000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| | | | | | | | Pyridine | 1000 ug/mL |
| ..SMLIST1 S10_00001 | 06/30/16 | | Restek, Lot A0107943 | | (Purchased Reagent) | | Benzoic acid | 2000 ug/mL |
| | | | | | | | Indene | 2000 ug/mL |
| ..SMLIST1 S11_00001 | 06/30/16 | | Restek, Lot A0108035 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1 S9_00001 | 07/31/16 | | Restek, Lot A0108174 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| | | | | | | | Benzidine | 2000 ug/mL |
| ..SMLIST1 SURR_00003 | 02/22/18 | | Restek, Lot A093638 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **SMLIST1 SS W_00005** | 04/20/16 | 04/21/15 | MECL2, Lot 0000093871 | 2 mL | SMIS80PPMW_00012 | 100 uL | 1,4-Dichlorobenzene-d4 | 4 ug/mL |
| | | | | | | | Acenaphthene-d10 | 4 ug/mL |
| | | | | | | | Chrysene-d12 | 4 ug/mL |
| | | | | | | | Naphthalene-d8 | 4 ug/mL |
| | | | | | | | Perylene-d12 | 4 ug/mL |
| | | | | | | | Phenanthrene-d10 | 4 ug/mL |
| .SMIS80PPMW_00012 | 04/20/16 | 04/20/15 | MECL2, Lot 0000073814 | 25 mL | SMIS R_00005 | 1 mL | 1,4-Dichlorobenzene-d4 | 80 ug/mL |
| | | | | | | | Acenaphthene-d10 | 80 ug/mL |
| | | | | | | | Chrysene-d12 | 80 ug/mL |
| | | | | | | | Naphthalene-d8 | 80 ug/mL |
| | | | | | | | Perylene-d12 | 80 ug/mL |
| | | | | | | | Phenanthrene-d10 | 80 ug/mL |
| ..SMIS R_00005 | 01/31/19 | | Restek, Lot A0101020 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2000 ug/mL |
| | | | | | | | Acenaphthene-d10 | 2000 ug/mL |
| | | | | | | | Chrysene-d12 | 2000 ug/mL |
| | | | | | | | Naphthalene-d8 | 2000 ug/mL |
| | | | | | | | Perylene-d12 | 2000 ug/mL |
| | | | | | | | Phenanthrene-d10 | 2000 ug/mL |
| **SMLIST1 SS W_00005** | 04/20/16 | 04/21/15 | MECL2, Lot 0000093871 | 2 mL | SMLIST1 SS ST_00005 | 200 uL | 1,1'-Biphenyl | 10 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 10 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 10 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 20 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 10 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 10 ug/mL |
| | | | | | | | 2-Chlorophenol | 10 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 10 ug/mL |
| | | | | | | | 2-Methylphenol | 10 ug/mL |
| | | | | | | | 2-Nitroaniline | 10 ug/mL |
| | | | | | | | 2-Nitrophenol | 10 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 10 ug/mL |
| | | | | | | | 3-Nitroaniline | 10 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 20 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 10 ug/mL |
| | | | | | | | 4-Chloroaniline | 10 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Chlorophenyl phenyl ether | 10 ug/mL |
| | | | | | | | 4-Nitroaniline | 10 ug/mL |
| | | | | | | | 4-Nitrophenol | 20 ug/mL |
| | | | | | | | Acenaphthene | 10 ug/mL |
| | | | | | | | Acenaphthylene | 10 ug/mL |
| | | | | | | | Acetophenone | 10 ug/mL |
| | | | | | | | Anthracene | 10 ug/mL |
| | | | | | | | Benzo[a]anthracene | 10 ug/mL |
| | | | | | | | Benzo[a]pyrene | 10 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 10 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 10 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 10 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 10 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 10 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 10 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 10 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 10 ug/mL |
| | | | | | | | Carbazole | 10 ug/mL |
| | | | | | | | Chrysene | 10 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 10 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 10 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 10 ug/mL |
| | | | | | | | Dibenzofuran | 10 ug/mL |
| | | | | | | | Diethyl phthalate | 10 ug/mL |
| | | | | | | | Dimethyl phthalate | 10 ug/mL |
| | | | | | | | Fluoranthene | 10 ug/mL |
| | | | | | | | Fluorene | 10 ug/mL |
| | | | | | | | Hexachlorobenzene | 10 ug/mL |
| | | | | | | | Hexachlorobutadiene | 10 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 10 ug/mL |
| | | | | | | | Hexachloroethane | 10 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 10 ug/mL |
| | | | | | | | Isophorone | 10 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 10 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 20 ug/mL |
| | | | | | | | Naphthalene | 10 ug/mL |
| | | | | | | | Nitrobenzene | 10 ug/mL |
| | | | | | | | Pentachlorophenol | 20 ug/mL |
| | | | | | | | Phenanthrene | 10 ug/mL |
| | | | | | | | Phenol | 10 ug/mL |
| | | | | | | | Pyrene | 10 ug/mL |
| | | | | | | | 3,3'-Dichlorobenzidine | 20 ug/mL |
| | | | | | | | Atrazine | 20 ug/mL |
| | | | | | | | Benzaldehyde | 20 ug/mL |
| | | | | | | | Caprolactam | 20 ug/mL |
| | | | | | | | 2,4,6-Tribromophenol (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 10 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 10 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Nitrobenzene-d5 (Surr) | 10 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 10 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 10 ug/mL |
| .SMLIST1 SS ST_00005 | 06/30/16 | 04/21/15 | MECL2, Lot 0000103549 | 10 mL | SMLIST1 SS S1_00004 | 1 mL | 1,1'-Biphenyl | 100 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 100 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 100 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 200 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 100 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 100 ug/mL |
| | | | | | | | 2-Chlorophenol | 100 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| | | | | | | | 2-Methylphenol | 100 ug/mL |
| | | | | | | | 2-Nitroaniline | 100 ug/mL |
| | | | | | | | 2-Nitrophenol | 100 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 100 ug/mL |
| | | | | | | | 3-Nitroaniline | 100 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 200 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 100 ug/mL |
| | | | | | | | 4-Chloroaniline | 100 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 100 ug/mL |
| | | | | | | | 4-Nitroaniline | 100 ug/mL |
| | | | | | | | 4-Nitrophenol | 200 ug/mL |
| | | | | | | | Acenaphthene | 100 ug/mL |
| | | | | | | | Acenaphthylene | 100 ug/mL |
| | | | | | | | Acetophenone | 100 ug/mL |
| | | | | | | | Anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]anthracene | 100 ug/mL |
| | | | | | | | Benzo[a]pyrene | 100 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 100 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 100 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 100 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 100 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 100 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 100 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 100 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 100 ug/mL |
| | | | | | | | Carbazole | 100 ug/mL |
| | | | | | | | Chrysene | 100 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 100 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 100 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 100 ug/mL |
| | | | | | | | Dibenzofuran | 100 ug/mL |
| | | | | | | | Diethyl phthalate | 100 ug/mL |
| | | | | | | | Dimethyl phthalate | 100 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Fluoranthene | 100 ug/mL |
| | | | | | | | Fluorene | 100 ug/mL |
| | | | | | | | Hexachlorobenzene | 100 ug/mL |
| | | | | | | | Hexachlorobutadiene | 100 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 100 ug/mL |
| | | | | | | | Hexachloroethane | 100 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 100 ug/mL |
| | | | | | | | Isophorone | 100 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 100 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 200 ug/mL |
| | | | | | | | Naphthalene | 100 ug/mL |
| | | | | | | | Nitrobenzene | 100 ug/mL |
| | | | | | | | Pentachlorophenol | 200 ug/mL |
| | | | | | | | Phenanthrene | 100 ug/mL |
| | | | | | | | Phenol | 100 ug/mL |
| | | | | | | | Pyrene | 100 ug/mL |
| | | | | | SMLIST1 SS S9_00001 | 1 mL | 3,3'-Dichlorobenzidine | 200 ug/mL |
| | | | | | SMLIST1SS S11_00001 | 1 mL | Atrazine | 200 ug/mL |
| | | | | | | | Benzaldehyde | 200 ug/mL |
| | | | | | | | Caprolactam | 200 ug/mL |
| | | | | | SMLIST1SS SUR_00002 | 200 uL | 2,4,6-Tribromophenol (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 100 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 100 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 100 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 100 ug/mL |
| | | | | | | | Terphenyl-d14 (Surr) | 100 ug/mL |
| ..SMLIST1 SS S1_00004 | 06/30/16 | | Restek, Lot A0107697 | | (Purchased Reagent) | | 1,1'-Biphenyl | 1000 ug/mL |
| | | | | | | | 2,4,5-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4,6-Trichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dichlorophenol | 1000 ug/mL |
| | | | | | | | 2,4-Dimethylphenol | 1000 ug/mL |
| | | | | | | | 2,4-Dinitrophenol | 2000 ug/mL |
| | | | | | | | 2,4-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2,6-Dinitrotoluene | 1000 ug/mL |
| | | | | | | | 2-Chloronaphthalene | 1000 ug/mL |
| | | | | | | | 2-Chlorophenol | 1000 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 1000 ug/mL |
| | | | | | | | 2-Methylphenol | 1000 ug/mL |
| | | | | | | | 2-Nitroaniline | 1000 ug/mL |
| | | | | | | | 2-Nitrophenol | 1000 ug/mL |
| | | | | | | | 3 & 4 Methylphenol | 1000 ug/mL |
| | | | | | | | 3-Nitroaniline | 1000 ug/mL |
| | | | | | | | 4,6-Dinitro-2-methylphenol | 2000 ug/mL |
| | | | | | | | 4-Bromophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Chloro-3-methylphenol | 1000 ug/mL |
| | | | | | | | 4-Chloroaniline | 1000 ug/mL |
| | | | | | | | 4-Chlorophenyl phenyl ether | 1000 ug/mL |
| | | | | | | | 4-Nitroaniline | 1000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 4-Nitrophenol | 2000 ug/mL |
| | | | | | | | Acenaphthene | 1000 ug/mL |
| | | | | | | | Acenaphthylene | 1000 ug/mL |
| | | | | | | | Acetophenone | 1000 ug/mL |
| | | | | | | | Anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]anthracene | 1000 ug/mL |
| | | | | | | | Benzo[a]pyrene | 1000 ug/mL |
| | | | | | | | Benzo[b]fluoranthene | 1000 ug/mL |
| | | | | | | | Benzo[g,h,i]perylene | 1000 ug/mL |
| | | | | | | | Benzo[k]fluoranthene | 1000 ug/mL |
| | | | | | | | bis (2-chloroisopropyl) ether | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethoxy)methane | 1000 ug/mL |
| | | | | | | | Bis(2-chloroethyl)ether | 1000 ug/mL |
| | | | | | | | Bis(2-ethylhexyl) phthalate | 1000 ug/mL |
| | | | | | | | Butyl benzyl phthalate | 1000 ug/mL |
| | | | | | | | Carbazole | 1000 ug/mL |
| | | | | | | | Chrysene | 1000 ug/mL |
| | | | | | | | Di-n-butyl phthalate | 1000 ug/mL |
| | | | | | | | Di-n-octyl phthalate | 1000 ug/mL |
| | | | | | | | Dibenz(a,h)anthracene | 1000 ug/mL |
| | | | | | | | Dibenzofuran | 1000 ug/mL |
| | | | | | | | Diethyl phthalate | 1000 ug/mL |
| | | | | | | | Dimethyl phthalate | 1000 ug/mL |
| | | | | | | | Fluoranthene | 1000 ug/mL |
| | | | | | | | Fluorene | 1000 ug/mL |
| | | | | | | | Hexachlorobenzene | 1000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 1000 ug/mL |
| | | | | | | | Hexachlorocyclopentadiene | 1000 ug/mL |
| | | | | | | | Hexachloroethane | 1000 ug/mL |
| | | | | | | | Indeno[1,2,3-cd]pyrene | 1000 ug/mL |
| | | | | | | | Isophorone | 1000 ug/mL |
| | | | | | | | N-Nitrosodi-n-propylamine | 1000 ug/mL |
| | | | | | | | N-Nitrosodiphenylamine | 2000 ug/mL |
| | | | | | | | Naphthalene | 1000 ug/mL |
| | | | | | | | Nitrobenzene | 1000 ug/mL |
| | | | | | | | Pentachlorophenol | 2000 ug/mL |
| | | | | | | | Phenanthrene | 1000 ug/mL |
| | | | | | | | Phenol | 1000 ug/mL |
| | | | | | | | Pyrene | 1000 ug/mL |
| ..SMLIST1 SS S9_00001 | 07/31/16 | | Restek, Lot A0108170 | | (Purchased Reagent) | | 3,3'-Dichlorobenzidine | 2000 ug/mL |
| ..SMLIST1SS S11_00001 | 06/30/16 | | Restek, Lot A0108042 | | (Purchased Reagent) | | Atrazine | 2000 ug/mL |
| | | | | | | | Benzaldehyde | 2000 ug/mL |
| | | | | | | | Caprolactam | 2000 ug/mL |
| ..SMLIST1SS SUR_00002 | 09/30/18 | | Restek, Lot A092712 | | (Purchased Reagent) | | 2,4,6-Tribromophenol (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorobiphenyl (Surr) | 5000 ug/mL |
| | | | | | | | 2-Fluorophenol (Surr) | 5000 ug/mL |
| | | | | | | | Nitrobenzene-d5 (Surr) | 5000 ug/mL |
| | | | | | | | Phenol-d5 (Surr) | 5000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Terphenyl-d14 (Surr) | 5000 ug/mL |
| **VGGASCALNEW_00015** | 12/23/15 | 09/23/15 | MEOH, Lot 62345 | 25 mL | VGGASSTOCKNEW_00003 | 1 mL | C6-C12 | 100 ug/mL |
| | | | | | | | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGGASSTOCKNEW_00003 | 01/31/20 | | RESTEK, Lot A092645 | | (Purchased Reagent) | | C6-C12 | 2500 ug/mL |
| | | | | | | | Gasoline Range Organics [C6 - C10] | 2500 ug/mL |
| **VGGASCALNEW_00016** | 03/08/16 | 12/08/15 | MEOH, Lot 62345 | 25 mL | VGGASSTOCKNEW_00003 | 1 mL | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGGASSTOCKNEW_00003 | 01/31/20 | | RESTEK, Lot A092645 | | (Purchased Reagent) | | Gasoline Range Organics [C6 - C10] | 2500 ug/mL |
| **VGGASCHK_00070** | 02/04/16 | 11/04/15 | MEOH, Lot 49909 | 50 mL | VGASCHK_00002 | 1 mL | Gasoline Range Organics [C6 - C10] | 100 ug/mL |
| .VGASCHK_00002 | 10/30/19 | | AccuStandard, Lot B5040009-01 | | (Purchased Reagent) | | Gasoline Range Organics [C6 - C10] | 5 mg/mL |
| **VGGASURROGATE_00034** | 01/09/16 | 10/09/15 | MEOH, Lot 49909 | 50 mL | VGSURRSTKGAS_00019 | 1.25 mL | Trifluorotoluene (Surr) | 50 ug/mL |
| .VGSURRSTKGAS_00019 | 08/31/16 | | ULTRA SCIENTIFIC, Lot CE-1903A | | (Purchased Reagent) | | Trifluorotoluene (Surr) | 2000 ug/mL |
| **VM250-IS_00031** | 04/21/16 | 11/21/15 | MEOH, Lot +171940000118655jm | 100 mL | vmcus8126_00008 | 1 mL | 1,4-Dichlorobenzene-d4 | 250 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 250 ug/mL |
| | | | | | | | Fluorobenzene | 250 ug/mL |
| .vmcus8126_00008 | 09/30/16 | | Ultra Scientific, Lot CL-3192 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 25000 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 25000 ug/mL |
| | | | | | | | Fluorobenzene | 25000 ug/mL |
| **VM250SS_00033** | 02/09/16 | 11/09/15 | MEOH, Lot +171940000118655JM | 50 mL | VM567650_00023 | 5 mL | 1,2-Dichloroethane-d4 (Surr) | 250 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 250 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 250 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 250 ug/mL |
| .VM567650_00023 | 06/30/19 | | Restek, Lot A0104073 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **VM50IS_00053** | 05/24/16 | 11/24/15 | MEOH, Lot 118655 | 100 mL | vm30241_00002 | 2 mL | 1,4-Dichlorobenzene-d4 | 50 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 50 ug/mL |
| | | | | | | | Fluorobenzene | 50 ug/mL |
| .vm30241_00002 | 11/30/19 | | restek, Lot A0107133 | | (Purchased Reagent) | | 1,4-Dichlorobenzene-d4 | 2500 ug/mL |
| | | | | | | | Chlorobenzene-d5 | 2500 ug/mL |
| | | | | | | | Fluorobenzene | 2500 ug/mL |
| **vm50ss_00225** | 01/04/16 | 12/28/15 | MEOH, Lot na | 2 mL | vm50ss_stk_00067 | 2 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| .vm50ss_stk_00067 | 04/05/16 | 10/05/15 | MEOH, Lot 0000099494 | 200 mL | VM567650_00023 | 4 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| ..VM567650_00023 | 06/30/19 | | Restek, Lot A0104073 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **vm50ss_stk_00067** | 04/05/16 | 10/05/15 | MEOH, Lot 0000099494 | 200 mL | VM567650_00023 | 4 mL | 1,2-Dichloroethane-d4 (Surr) | 50 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 50 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 50 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 50 ug/mL |
| .VM567650_00023 | 06/30/19 | | Restek, Lot A0104073 | | (Purchased Reagent) | | 1,2-Dichloroethane-d4 (Surr) | 2500 ug/mL |
| | | | | | | | 4-Bromofluorobenzene (Surr) | 2500 ug/mL |
| | | | | | | | Dibromofluoromethane (Surr) | 2500 ug/mL |
| | | | | | | | Toluene-d8 (Surr) | 2500 ug/mL |
| **VMAROLISTDW_00128** | 12/16/15 | 12/09/15 | MEOH, Lot na | 3 mL | VMACROLSTD_00032 | 3 mL | Acrolein | 250 ug/mL |
| .VMACROLSTD_00032 | 02/29/16 | 11/30/15 | MEOH, Lot +171940000118655jm | 20 mL | VM568720_00010 | 250 uL | Acrolein | 250 ug/mL |
| ..VM568720_00010 | 03/31/16 | | restek, Lot A0115119 | | (Purchased Reagent) | | Acrolein | 20000 ug/mL |
| **VMAROLISTDW_00131** | 01/04/16 | 12/28/15 | MEOH, Lot na | 3 mL | VMACROLSTD_00032 | 3 mL | Acrolein | 250 ug/mL |
| .VMACROLSTD_00032 | 02/29/16 | 11/30/15 | MEOH, Lot +171940000118655jm | 20 mL | VM568720_00010 | 250 uL | Acrolein | 250 ug/mL |
| ..VM568720_00010 | 03/31/16 | | restek, Lot A0115119 | | (Purchased Reagent) | | Acrolein | 20000 ug/mL |
| **VMFASA9W_00080** | 12/16/15 | 12/09/15 | MEOH, Lot NA | 2 mL | VMFASA9_00006 | 2 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| .VMFASA9_00006 | 02/29/16 | 11/30/15 | MEOH, Lot +171940000118655jm | 100 mL | vm569725S_00003 | 2 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | VM569728S_00001 | 2 mL | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| ..vm569725S_00003 | 02/29/16 | | restek, Lot A0113467 | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| ..VM569728S_00001 | 04/30/17 | | restek, Lot A0110751 | | (Purchased Reagent) | | Diisopropyl ether | 2500 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2500 ug/mL |
| **VMFASA9W_00082** | 01/05/16 | 12/29/15 | MEOH, Lot NA | 2 mL | VMFASA9_00006 | 2 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| .VMFASA9_00006 | 02/29/16 | 11/30/15 | MEOH, Lot +171940000118655jm | 100 mL | vm569725S_00003 | 2 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | VM569728S_00001 | 2 mL | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| ..vm569725S_00003 | 02/29/16 | | restek, Lot A0113467 | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| ..VM569728S_00001 | 04/30/17 | | restek, Lot A0110751 | | (Purchased Reagent) | | Diisopropyl ether | 2500 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **VMFASGW_00129** | 12/18/15 | 12/11/15 | MEOH, Lot NA | 2 mL | VMFASG_00037 | 2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .VMFASG_00037 | 12/30/15 | 11/30/15 | MEOH, Lot +171940000118655jm | 50 mL | vm569722S_00001 | 1 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| ..vm569722S_00001 | 08/31/18 | | Restek, Lot A0113063 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMFASGW_00132** | 01/06/16 | 12/30/15 | MEOH, Lot NA | 2 mL | VMFASG_00038 | 2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .VMFASG_00038 | 01/31/16 | 12/30/15 | MEOH, Lot 0000118655 | 50 mL | vm569722S_00001 | 1 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| ..vm569722S_00001 | 08/31/18 | | Restek, Lot A0113063 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMFASPW_00130** | 12/18/15 | 12/11/15 | MEOH, Lot n/a | 2 mL | VMRFASP_00024 | 2 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRFASP_00024 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 100 mL | VM567641S_00007 | 2.5 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | VM567642S_00005 | 1 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM569723S_00001 | 2 mL | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | VM569724S_00003 | 1 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641S_00007 | 02/29/16 | | Restek, Lot A093733 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| | | | | | | | Xylenes, Total | 4000 ug/mL |
| ..VM567642S_00005 | 02/29/16 | | Restek, Lot A093472 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM569723S_00001 | 01/31/18 | | Restek, Lot A0108188 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2500 ug/mL |
| ..VM569724S_00003 | 01/31/16 | | Restek, Lot A0112364 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMFASPW_00133** | 01/06/16 | 12/30/15 | MEOH, Lot n/a | 2 mL | VMRFASP_00025 | 2 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRFASP_00025 | 01/30/16 | 12/30/15 | MEOH, Lot +171940000118655JM | 100 mL | VM567642S_00005 | 1 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM569720S_00001 | 2 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | Xylenes, Total | 100 ug/mL |
| | | | | | VM569723S_00001 | 2 mL | 2-Chloroethyl vinyl ether | 50 ug/mL |
| | | | | | VM569724S_00003 | 1 mL | Vinyl acetate | 50 ug/mL |
| ..VM567642S_00005 | 02/29/16 | | Restek, Lot A093472 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM569720S_00001 | 01/31/17 | | Restek, Lot A0108163 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2500 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2500 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2500 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2500 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2500 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2500 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2500 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2500 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2500 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2500 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2500 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2500 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2500 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2500 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2500 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2500 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2500 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2500 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2500 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2500 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2500 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2500 ug/mL |
| | | | | | | | Benzene | 2500 ug/mL |
| | | | | | | | Bromobenzene | 2500 ug/mL |
| | | | | | | | Bromochloromethane | 2500 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Bromodichloromethane | 2500 ug/mL |
| | | | | | | | Bromoform | 2500 ug/mL |
| | | | | | | | Carbon disulfide | 2500 ug/mL |
| | | | | | | | Carbon tetrachloride | 2500 ug/mL |
| | | | | | | | Chlorobenzene | 2500 ug/mL |
| | | | | | | | Chloroform | 2500 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2500 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2500 ug/mL |
| | | | | | | | Dibromochloromethane | 2500 ug/mL |
| | | | | | | | Dibromomethane | 2500 ug/mL |
| | | | | | | | Ethylbenzene | 2500 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2500 ug/mL |
| | | | | | | | Isopropylbenzene | 2500 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2500 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2500 ug/mL |
| | | | | | | | Methylene Chloride | 2500 ug/mL |
| | | | | | | | n-Butylbenzene | 2500 ug/mL |
| | | | | | | | N-Propylbenzene | 2500 ug/mL |
| | | | | | | | Naphthalene | 2500 ug/mL |
| | | | | | | | o-Xylene | 2500 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2500 ug/mL |
| | | | | | | | sec-Butylbenzene | 2500 ug/mL |
| | | | | | | | Styrene | 2500 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 25000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2500 ug/mL |
| | | | | | | | Tetrachloroethene | 2500 ug/mL |
| | | | | | | | Toluene | 2500 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2500 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2500 ug/mL |
| | | | | | | | Trichloroethene | 2500 ug/mL |
| | | | | | | | Xylenes, Total | 5000 ug/mL |
| ..VM569723S_00001 | 01/31/18 | | Restek, Lot A0108188 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2500 ug/mL |
| ..VM569724S_00003 | 01/31/16 | | Restek, Lot A0112364 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRA9W_00135** | 12/16/15 | 12/09/15 | MEOH, Lot NA | 1 mL | VMRA9_00014 | 1 mL | Cyclohexanone | 500 ug/mL |
| | | | | | | | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| .VMRA9_00014 | 12/31/15 | 12/01/15 | MEOH, Lot 00000118655 | 50 mL | VM567648_00014 | 1.25 mL | Cyclohexanone | 500 ug/mL |
| | | | | | VM568723_00001 | 1.25 mL | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725_00004 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | vm569726_00002 | 2 mL | 2-Methylnaphthalene | 100 ug/mL |
| ..VM567648_00014 | 02/29/16 | Restek, Lot A093361 | | | (Purchased Reagent) | | Cyclohexanone | 20000 ug/mL |
| ..VM568723_00001 | 12/31/15 | restek, Lot A099930 | | | (Purchased Reagent) | | Acetonitrile | 20000 ug/mL |
| | | | | | | | Diisopropyl ether | 2000 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Propionitrile | 20000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725_00004 | 03/31/16 | restek, Lot A0113771 | | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 2500 ug/mL |
| | | | | | | | 1-Chlorohexane | 2500 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 2500 ug/mL |
| | | | | | | | 2-Nitropropane | 5000 ug/mL |
| | | | | | | | Ethyl acetate | 5000 ug/mL |
| | | | | | | | Ethyl acrylate | 2500 ug/mL |
| | | | | | | | Methacrylonitrile | 25000 ug/mL |
| | | | | | | | Methyl methacrylate | 5000 ug/mL |
| | | | | | | | n-Butanol | 62500 ug/mL |
| | | | | | | | n-Butyl acetate | 2500 ug/mL |
| ..vm569726_00002 | 01/31/17 | Restek, Lot A0108210 | | | (Purchased Reagent) | | 2-Methylnaphthalene | 2500 ug/mL |
| **VMRA9W_00137** | 12/31/15 | 12/28/15 | MEOH, Lot NA | 1 mL | VMRA9_00014 | 1 mL | Cyclohexanone | 500 ug/mL |
| | | | | | | | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | | | 1,2,3-Trimethylbenzene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | | | 2-Methylnaphthalene | 100 ug/mL |
| .VMRA9_00014 | 12/31/15 | 12/01/15 | MEOH, Lot 00000118655 | 50 mL | VM567648_00014 | 1.25 mL | Cyclohexanone | 500 ug/mL |
| | | | | | VM568723_00001 | 1.25 mL | Acetonitrile | 500 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Propionitrile | 500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| | | | | | vm569725_00004 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1-Chlorohexane | 50 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 50 ug/mL |
| | | | | | | | 2-Nitropropane | 100 ug/mL |
| | | | | | | | Ethyl acetate | 100 ug/mL |
| | | | | | | | Ethyl acrylate | 50 ug/mL |
| | | | | | | | Methacrylonitrile | 500 ug/mL |
| | | | | | | | Methyl methacrylate | 100 ug/mL |
| | | | | | | | n-Butanol | 1250 ug/mL |
| | | | | | | | n-Butyl acetate | 50 ug/mL |
| | | | | | vm569726_00002 | 2 mL | 2-Methylnaphthalene | 100 ug/mL |
| ..VM567648_00014 | 02/29/16 | Restek, Lot A093361 | | | (Purchased Reagent) | | Cyclohexanone | 20000 ug/mL |
| ..VM568723_00001 | 12/31/15 | restek, Lot A099930 | | | (Purchased Reagent) | | Acetonitrile | 20000 ug/mL |
| | | | | | | | Diisopropyl ether | 2000 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2000 ug/mL |
| | | | | | | | Propionitrile | 20000 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2000 ug/mL |
| ..vm569725_00004 | 03/31/16 | restek, Lot A0113771 | | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| | | | | | | | 1,3,5-Trichlorobenzene | 2500 ug/mL |
| | | | | | | | 1-Chlorohexane | 2500 ug/mL |
| | | | | | | | 2-Chloro-1,3-butadiene | 2500 ug/mL |
| | | | | | | | 2-Nitropropane | 5000 ug/mL |
| | | | | | | | Ethyl acetate | 5000 ug/mL |
| | | | | | | | Ethyl acrylate | 2500 ug/mL |
| | | | | | | | Methacrylonitrile | 25000 ug/mL |
| | | | | | | | Methyl methacrylate | 5000 ug/mL |
| | | | | | | | n-Butanol | 62500 ug/mL |
| | | | | | | | n-Butyl acetate | 2500 ug/mL |
| ..vm569726_00002 | 01/31/17 | Restek, Lot A0108210 | | | (Purchased Reagent) | | 2-Methylnaphthalene | 2500 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| **VMRA9W_00138** | 01/07/16 | 12/31/15 | MEOH, Lot NA | 1 mL | VMRA9_00015 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | | | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| .VMRA9_00015 | 03/31/16 | 12/31/15 | MEOH, Lot 00000118655 | 50 mL | vm569725_00004 | 1 mL | 1,2,3-Trimethylbenzene | 50 ug/mL |
| | | | | | VM569728_00001 | 1 mL | Diisopropyl ether | 50 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 50 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 50 ug/mL |
| ..vm569725_00004 | 03/31/16 | | restek, Lot A0113771 | | (Purchased Reagent) | | 1,2,3-Trimethylbenzene | 2500 ug/mL |
| ..VM569728_00001 | 04/30/17 | | restek, Lot A0110540 | | (Purchased Reagent) | | Diisopropyl ether | 2500 ug/mL |
| | | | | | | | Ethyl t-butyl ether | 2500 ug/mL |
| | | | | | | | Tert-amyl methyl ether | 2500 ug/mL |
| **VMRGAS_00129** | 12/15/15 | 12/08/15 | MEOH, Lot 0000118655 | 10 mL | vm569722_00003 | 0.2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Butadiene | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Dichlorofluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .vm569722_00003 | 04/30/18 | | Restek, Lot A0110070 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Butadiene | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Dichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMRGAS_00131** | 12/31/15 | 12/24/15 | MEOH, Lot 0000118655 | 10 mL | vm569722_00003 | 0.2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Butadiene | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Dichlorofluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .vm569722_00003 | 04/30/18 | | Restek, Lot A0110070 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Butadiene | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Dichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMRGAS_00132** | 01/09/16 | 01/02/16 | MEOH, Lot 0000118655 | 10 mL | vm569722_00004 | 0.2 mL | Bromomethane | 50 ug/mL |
| | | | | | | | Chloroethane | 50 ug/mL |
| | | | | | | | Chloromethane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Dichlorodifluoromethane | 50 ug/mL |
| | | | | | | | Trichlorofluoromethane | 50 ug/mL |
| | | | | | | | Vinyl chloride | 50 ug/mL |
| .vm569722_00004 | 04/30/18 | | Restek, Lot A0110070 | | (Purchased Reagent) | | Bromomethane | 2500 ug/mL |
| | | | | | | | Chloroethane | 2500 ug/mL |
| | | | | | | | Chloromethane | 2500 ug/mL |
| | | | | | | | Dichlorodifluoromethane | 2500 ug/mL |
| | | | | | | | Trichlorofluoromethane | 2500 ug/mL |
| | | | | | | | Vinyl chloride | 2500 ug/mL |
| **VMRPRIMW_00160** | 12/14/15 | 12/07/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00014 | 1 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRPRIM_00014 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 50 mL | VM567641_00013 | 1.25 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | VM567642_00015 | 0.5 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM567643_00020 | 2.5 mL | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | VM569724_00003 | 0.5 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641_00013 | 02/29/16 | | restek, Lot A093581 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,4-Dioxane | 40000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acrylonitrile | 20000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |
| | | | | | | | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Cyclohexane | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethyl ether | 2000 ug/mL |
| | | | | | | | Ethyl methacrylate | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Hexane | 2000 ug/mL |
| | | | | | | | Iodomethane | 2000 ug/mL |
| | | | | | | | Isobutyl alcohol | 50000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl acetate | 10000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylcyclohexane | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | n-Heptane | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Tetrahydrofuran | 4000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| ..VM567642_00015 | 02/29/16 | | Restek, Lot A093365 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM567643_00020 | 02/29/16 | | restek, Lot A093368 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2000 ug/mL |
| ..VM569724_00003 | 01/31/16 | | Restek, Lot A0112975 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRPRIMW_00163** | 01/05/16 | 12/29/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00014 | 1 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | | | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | | | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | | | Vinyl acetate | 50 ug/mL |
| .VMRPRIM_00014 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 50 mL | VM567641_00013 | 1.25 mL | 1,1,1,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 50 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 50 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 50 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 50 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | 1,2-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 50 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 50 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 50 ug/mL |
| | | | | | | | 1,4-Dioxane | 1000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 50 ug/mL |
| | | | | | | | 2-Chlorotoluene | 50 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 50 ug/mL |
| | | | | | | | 4-Chlorotoluene | 50 ug/mL |
| | | | | | | | Acrylonitrile | 500 ug/mL |
| | | | | | | | Benzene | 50 ug/mL |
| | | | | | | | Bromobenzene | 50 ug/mL |
| | | | | | | | Bromochloromethane | 50 ug/mL |
| | | | | | | | Bromodichloromethane | 50 ug/mL |
| | | | | | | | Bromoform | 50 ug/mL |
| | | | | | | | Carbon disulfide | 50 ug/mL |
| | | | | | | | Carbon tetrachloride | 50 ug/mL |
| | | | | | | | Chlorobenzene | 50 ug/mL |
| | | | | | | | Chloroform | 50 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | Cyclohexane | 50 ug/mL |
| | | | | | | | Dibromochloromethane | 50 ug/mL |
| | | | | | | | Dibromomethane | 50 ug/mL |
| | | | | | | | Ethyl ether | 50 ug/mL |
| | | | | | | | Ethyl methacrylate | 50 ug/mL |
| | | | | | | | Ethylbenzene | 50 ug/mL |
| | | | | | | | Hexachlorobutadiene | 50 ug/mL |
| | | | | | | | Hexane | 50 ug/mL |
| | | | | | | | Iodomethane | 50 ug/mL |
| | | | | | | | Isobutyl alcohol | 1250 ug/mL |
| | | | | | | | Isopropylbenzene | 50 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 50 ug/mL |
| | | | | | | | Methyl acetate | 250 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 50 ug/mL |
| | | | | | | | Methylcyclohexane | 50 ug/mL |
| | | | | | | | Methylene Chloride | 50 ug/mL |
| | | | | | | | n-Butylbenzene | 50 ug/mL |
| | | | | | | | n-Heptane | 50 ug/mL |
| | | | | | | | N-Propylbenzene | 50 ug/mL |
| | | | | | | | Naphthalene | 50 ug/mL |
| | | | | | | | o-Xylene | 50 ug/mL |
| | | | | | | | p-Isopropyltoluene | 50 ug/mL |
| | | | | | | | sec-Butylbenzene | 50 ug/mL |
| | | | | | | | Styrene | 50 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | | |
| | | | | | | | tert-Butyl alcohol (TBA) | 500 ug/mL |
| | | | | | | | tert-Butylbenzene | 50 ug/mL |
| | | | | | | | Tetrachloroethene | 50 ug/mL |
| | | | | | | | Tetrahydrofuran | 100 ug/mL |
| | | | | | | | Toluene | 50 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 50 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 50 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 50 ug/mL |
| | | | | | | | Trichloroethene | 50 ug/mL |
| | | | | | VM567642_00015 | 0.5 mL | 2-Butanone (MEK) | 100 ug/mL |
| | | | | | | | 2-Hexanone | 100 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 100 ug/mL |
| | | | | | | | Acetone | 100 ug/mL |
| | | | | | VM567643_00020 | 2.5 mL | 2-Chloroethyl vinyl ether | 100 ug/mL |
| | | | | | VM569724_00003 | 0.5 mL | Vinyl acetate | 50 ug/mL |
| ..VM567641_00013 | 02/29/16 | | restek, Lot A093581 | | (Purchased Reagent) | | 1,1,1,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,1-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2,2-Tetrachloroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloro-1,2,2-trifluoroethane | 2000 ug/mL |
| | | | | | | | 1,1,2-Trichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,1-Dichloroethene | 2000 ug/mL |
| | | | | | | | 1,1-Dichloropropene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,3-Trichloropropane | 2000 ug/mL |
| | | | | | | | 1,2,4-Trichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2,4-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dibromo-3-Chloropropane | 2000 ug/mL |
| | | | | | | | 1,2-Dibromoethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,2-Dichloroethane | 2000 ug/mL |
| | | | | | | | 1,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,3,5-Trimethylbenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,3-Dichloropropane | 2000 ug/mL |
| | | | | | | | 1,4-Dichlorobenzene | 2000 ug/mL |
| | | | | | | | 1,4-Dioxane | 40000 ug/mL |
| | | | | | | | 2,2-Dichloropropane | 2000 ug/mL |
| | | | | | | | 2-Chlorotoluene | 2000 ug/mL |
| | | | | | | | 3-Chloro-1-propene | 2000 ug/mL |
| | | | | | | | 4-Chlorotoluene | 2000 ug/mL |
| | | | | | | | Acrylonitrile | 20000 ug/mL |
| | | | | | | | Benzene | 2000 ug/mL |
| | | | | | | | Bromobenzene | 2000 ug/mL |
| | | | | | | | Bromochloromethane | 2000 ug/mL |
| | | | | | | | Bromodichloromethane | 2000 ug/mL |
| | | | | | | | Bromoform | 2000 ug/mL |

# REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

| Reagent ID | Exp Date | Prep Date | Dilutant Used | Reagent Final Volume | Parent Reagent | | Analyte | Concentration |
|---|---|---|---|---|---|---|---|---|
| | | | | | Reagent ID | Volume Added | Carbon disulfide | 2000 ug/mL |
| | | | | | | | Carbon tetrachloride | 2000 ug/mL |
| | | | | | | | Chlorobenzene | 2000 ug/mL |
| | | | | | | | Chloroform | 2000 ug/mL |
| | | | | | | | cis-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | cis-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | Cyclohexane | 2000 ug/mL |
| | | | | | | | Dibromochloromethane | 2000 ug/mL |
| | | | | | | | Dibromomethane | 2000 ug/mL |
| | | | | | | | Ethyl ether | 2000 ug/mL |
| | | | | | | | Ethyl methacrylate | 2000 ug/mL |
| | | | | | | | Ethylbenzene | 2000 ug/mL |
| | | | | | | | Hexachlorobutadiene | 2000 ug/mL |
| | | | | | | | Hexane | 2000 ug/mL |
| | | | | | | | Iodomethane | 2000 ug/mL |
| | | | | | | | Isobutyl alcohol | 50000 ug/mL |
| | | | | | | | Isopropylbenzene | 2000 ug/mL |
| | | | | | | | m-Xylene & p-Xylene | 2000 ug/mL |
| | | | | | | | Methyl acetate | 10000 ug/mL |
| | | | | | | | Methyl tert-butyl ether | 2000 ug/mL |
| | | | | | | | Methylcyclohexane | 2000 ug/mL |
| | | | | | | | Methylene Chloride | 2000 ug/mL |
| | | | | | | | n-Butylbenzene | 2000 ug/mL |
| | | | | | | | n-Heptane | 2000 ug/mL |
| | | | | | | | N-Propylbenzene | 2000 ug/mL |
| | | | | | | | Naphthalene | 2000 ug/mL |
| | | | | | | | o-Xylene | 2000 ug/mL |
| | | | | | | | p-Isopropyltoluene | 2000 ug/mL |
| | | | | | | | sec-Butylbenzene | 2000 ug/mL |
| | | | | | | | Styrene | 2000 ug/mL |
| | | | | | | | tert-Butyl alcohol (TBA) | 20000 ug/mL |
| | | | | | | | tert-Butylbenzene | 2000 ug/mL |
| | | | | | | | Tetrachloroethene | 2000 ug/mL |
| | | | | | | | Tetrahydrofuran | 4000 ug/mL |
| | | | | | | | Toluene | 2000 ug/mL |
| | | | | | | | trans-1,2-Dichloroethene | 2000 ug/mL |
| | | | | | | | trans-1,3-Dichloropropene | 2000 ug/mL |
| | | | | | | | trans-1,4-Dichloro-2-butene | 2000 ug/mL |
| | | | | | | | Trichloroethene | 2000 ug/mL |
| ..VM567642_00015 | 02/29/16 | | Restek, Lot A093365 | | (Purchased Reagent) | | 2-Butanone (MEK) | 10000 ug/mL |
| | | | | | | | 2-Hexanone | 10000 ug/mL |
| | | | | | | | 4-Methyl-2-pentanone (MIBK) | 10000 ug/mL |
| | | | | | | | Acetone | 10000 ug/mL |
| ..VM567643_00020 | 02/29/16 | | restek, Lot A093368 | | (Purchased Reagent) | | 2-Chloroethyl vinyl ether | 2000 ug/mL |
| ..VM569724_00003 | 01/31/16 | | Restek, Lot A0112975 | | (Purchased Reagent) | | Vinyl acetate | 5000 ug/mL |
| **VMRPRIMW_00163** | 01/05/16 | 12/29/15 | MEOH, Lot NA | 1 mL | VMRPRIM_00014 | 1 mL | Xylenes, Total | 100 ug/mL |
| .VMRPRIM_00014 | 01/31/16 | 11/01/15 | MEOH, Lot 0000118655 | 50 mL | VM567641_00013 | 1.25 mL | Xylenes, Total | 100 ug/mL |
| ..VM567641_00013 | 02/29/16 | | restek, Lot A093581 | | (Purchased Reagent) | | Xylenes, Total | 4000 ug/mL |

# Certification Summary

Client: Tetra Tech, Inc. TestAmerica Job ID: 240-59482-1
Project/Site: MRC Block F UST Pipe Inv.

| Laboratory | Authority | Program | EPA Region | Certification ID |
|---|---|---|---|---|
| TestAmerica Canton | California | State Program | 9 | 2927 |
| TestAmerica Canton | Connecticut | State Program | 1 | PH-0590 |
| TestAmerica Canton | Illinois | NELAP | 5 | 200004 |
| TestAmerica Canton | Kansas | NELAP | 7 | E-10336 |
| TestAmerica Canton | Kentucky (UST) | State Program | 4 | 58 |
| TestAmerica Canton | Kentucky (WW) | State Program | 4 | 98016 |
| TestAmerica Canton | L-A-B | DoD ELAP | | L2315 |
| TestAmerica Canton | Minnesota | NELAP | 5 | 039-999-348 |
| TestAmerica Canton | Nevada | State Program | 9 | OH-000482008A |
| TestAmerica Canton | New Jersey | NELAP | 2 | OH001 |
| TestAmerica Canton | New York | NELAP | 2 | 10975 |
| TestAmerica Canton | Ohio VAP | State Program | 5 | CL0024 |
| TestAmerica Canton | Oregon | NELAP | 10 | 4062 |
| TestAmerica Canton | Pennsylvania | NELAP | 3 | 68-00340 |
| TestAmerica Canton | Texas | NELAP | 6 | T104704517-15-5 |
| TestAmerica Canton | USDA | Federal | | P330-13-00319 |
| TestAmerica Canton | Virginia | NELAP | 3 | 460175 |
| TestAmerica Canton | Washington | State Program | 10 | C971 |
| TestAmerica Canton | West Virginia DEP | State Program | 3 | 210 |
| TestAmerica Canton | Wisconsin | State Program | 5 | 999518190 |

Accreditation may not be offered or required for all methods and analytes reported in this package. Please contact your project manager for the laboratory's current list of certified methods and analytes.

# Method 8260C

## Volatile Organic Compounds (GC/MS) by Method 8260C

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): DB-624 ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | DBFM # | DCA # | TOL # | BFB # |
|---|---|---|---|---|---|
| F-VB-UST-P-7 | 240-59482-2 | 93 | 100 | 91 | 88 |
| F-VB-UST-P-8 | 240-59482-3 | 90 | 97 | 91 | 87 |
| | MB 240-212606/24 | 93 | 97 | 88 | 86 |
| | LCS 240-212606/23 | 94 | 92 | 97 | 91 |
| | LCS 240-212606/25 | 90 | 92 | 93 | 89 |
| F-VB-UST-P-7 MS | 240-59482-2 MS | 95 | 97 | 92 | 91 |
| F-VB-UST-P-7 MSD | 240-59482-2 MSD | 95 | 100 | 91 | 87 |

| | QC LIMITS |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 37-132 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 58-123 |
| TOL = Toluene-d8 (Surr) | 67-125 |
| BFB = 4-Bromofluorobenzene (Surr) | 52-136 |

# Column to be used to flag recovery values

FORM II 8260C

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Water Level: Low

GC Column (1): DB-624 ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | DBFM # | DCA # | TOL # | BFB # |
|---|---|---|---|---|---|
| TB-122815 | 240-59482-1 | 93 | 99 | 102 | 112 |
| | MB 240-212893/8 | 94 | 97 | 102 | 111 |
| | LCS 240-212893/5 | 101 | 96 | 100 | 114 |
| | 240-59545-B-2 MS | 100 | 102 | 103 | 117 |
| | 240-59545-B-2 MSD | 105 | 97 | 98 | 116 |

| | QC LIMITS |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 79-120 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 78-125 |
| TOL = Toluene-d8 (Surr) | 80-120 |
| BFB = 4-Bromofluorobenzene (Surr) | 61-120 |

# Column to be used to flag recovery values

FORM II 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 147320.D

Lab ID: LCS 240-212606/23 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 50.0 | 58.6 | 117 | 80-120 | |
| 1,1,1-Trichloroethane | 50.0 | 54.0 | 108 | 77-126 | |
| 1,1,2,2-Tetrachloroethane | 50.0 | 53.5 | 107 | 77-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 50.0 | 45.3 | 91 | 80-138 | |
| 1,1-Dichloroethane | 50.0 | 51.1 | 102 | 76-120 | |
| 1,1-Dichloroethene | 50.0 | 42.2 | 84 | 75-135 | |
| 1,1-Dichloropropene | 50.0 | 54.6 | 109 | 80-120 | |
| 1,2,3-Trichlorobenzene | 50.0 | 45.6 | 91 | 61-121 | |
| 1,2,3-Trichloropropane | 50.0 | 57.2 | 114 | 80-126 | |
| 1,2,4-Trichlorobenzene | 50.0 | 45.6 | 91 | 64-124 | |
| 1,2,4-Trimethylbenzene | 50.0 | 52.6 | 105 | 80-129 | |
| 1,2-Dibromo-3-Chloropropane | 50.0 | 55.6 | 111 | 61-132 | |
| 1,2-Dichlorobenzene | 50.0 | 48.1 | 96 | 76-120 | |
| 1,2-Dichloroethane | 50.0 | 52.5 | 105 | 72-120 | |
| 1,2-Dichloropropane | 50.0 | 52.7 | 105 | 80-120 | |
| 1,3-Dichlorobenzene | 50.0 | 47.0 | 94 | 78-120 | |
| 1,3-Dichloropropane | 50.0 | 51.7 | 103 | 80-120 | |
| 1,4-Dichlorobenzene | 50.0 | 47.3 | 95 | 75-120 | |
| 2,2-Dichloropropane | 50.0 | 54.3 | 109 | 69-135 | |
| 2-Butanone (MEK) | 100 | 96.3 | 96 | 52-131 | |
| 2-Chloroethyl vinyl ether | 50.0 | 54.3 | 109 | 28-173 | |
| 2-Chlorotoluene | 50.0 | 52.2 | 104 | 78-120 | |
| 2-Hexanone | 100 | 99.2 | 99 | 64-136 | |
| 4-Chlorotoluene | 50.0 | 48.3 | 97 | 79-120 | |
| p-Isopropyltoluene | 50.0 | 53.4 | 107 | 77-131 | |
| 4-Methyl-2-pentanone (MIBK) | 100 | 109 | 109 | 67-135 | |
| Acetone | 100 | 94.9 | 95 | 41-137 | |
| Benzene | 50.0 | 50.8 | 102 | 79-120 | |
| Bromobenzene | 50.0 | 51.6 | 103 | 80-120 | |
| Bromoform | 50.0 | 54.9 | 110 | 62-133 | |
| Bromomethane | 20.0 | 19.2 | 96 | 42-136 | |
| Carbon disulfide | 50.0 | 54.9 | 110 | 62-146 | |
| Carbon tetrachloride | 50.0 | 57.7 | 115 | 71-129 | |
| Chlorobenzene | 50.0 | 50.5 | 101 | 78-120 | |
| Dibromochloromethane | 50.0 | 54.7 | 109 | 72-127 | |
| Chloroethane | 20.0 | 19.5 | 98 | 58-120 | |
| Chloroform | 50.0 | 51.8 | 104 | 77-120 | |
| Chloromethane | 20.0 | 20.5 | 102 | 50-120 | |
| cis-1,2-Dichloroethene | 50.0 | 52.6 | 105 | 76-120 | |
| cis-1,3-Dichloropropene | 50.0 | 53.5 | 107 | 74-128 | |
| Dibromomethane | 50.0 | 53.8 | 108 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 147320.D

Lab ID: LCS 240-212606/23 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 50.0 | 52.0 | 104 | 79-120 | |
| Dichlorodifluoromethane | 20.0 | 20.7 | 103 | 26-120 | |
| Bromodichloromethane | 50.0 | 54.4 | 109 | 80-122 | |
| tert-Butyl alcohol (TBA) | 500 | 450 | 90 | 60-125 | |
| Ethylbenzene | 50.0 | 51.6 | 103 | 79-120 | |
| Hexachlorobutadiene | 50.0 | 51.3 | 103 | 54-131 | |
| Isopropylbenzene | 50.0 | 54.4 | 109 | 76-122 | |
| Methyl tert-butyl ether | 50.0 | 53.0 | 106 | 49-165 | |
| Methylene Chloride | 50.0 | 50.7 | 101 | 75-120 | |
| m-Xylene & p-Xylene | 50.0 | 53.8 | 108 | 80-120 | |
| Naphthalene | 50.0 | 51.1 | 102 | 65-123 | |
| n-Butylbenzene | 50.0 | 50.3 | 101 | 68-135 | |
| N-Propylbenzene | 50.0 | 53.2 | 106 | 80-129 | |
| o-Xylene | 50.0 | 54.6 | 109 | 80-120 | |
| sec-Butylbenzene | 50.0 | 55.3 | 111 | 74-129 | |
| Styrene | 50.0 | 53.8 | 108 | 80-120 | |
| Tetrachloroethene | 50.0 | 51.4 | 103 | 79-120 | |
| Toluene | 50.0 | 53.2 | 106 | 75-120 | |
| trans-1,2-Dichloroethene | 50.0 | 53.2 | 106 | 78-120 | |
| trans-1,3-Dichloropropene | 50.0 | 55.5 | 111 | 73-131 | |
| Trichloroethene | 50.0 | 53.1 | 106 | 79-120 | |
| Trichlorofluoromethane | 20.0 | 21.1 | 106 | 57-146 | |
| Vinyl acetate | 50.0 | 48.6 | 97 | 70-130 | |
| Vinyl chloride | 20.0 | 20.2 | 101 | 57-120 | |
| Xylenes, Total | 100 | 108 | 108 | 80-120 | |
| tert-Butylbenzene | 50.0 | 54.9 | 110 | 76-126 | |
| 1,2-Dibromoethane | 50.0 | 56.1 | 112 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC5593.D

Lab ID: LCS 240-212893/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 10.0 | 8.23 | 82 | 80-120 | |
| 1,1,1-Trichloroethane | 10.0 | 8.79 | 88 | 77-123 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 9.40 | 94 | 71-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 10.0 | 10.5 | 105 | 67-138 | |
| 1,1-Dichloroethane | 10.0 | 9.36 | 94 | 79-125 | |
| 1,1-Dichloroethene | 10.0 | 9.64 | 96 | 76-124 | |
| 1,1-Dichloropropene | 10.0 | 10.1 | 101 | 80-128 | |
| 1,2,3-Trichlorobenzene | 10.0 | 8.36 | 84 | 66-120 | |
| 1,2,3-Trichloropropane | 10.0 | 9.73 | 97 | 72-126 | |
| 1,2,4-Trichlorobenzene | 10.0 | 8.53 | 85 | 61-120 | |
| 1,2,4-Trimethylbenzene | 10.0 | 9.27 | 93 | 76-120 | |
| 1,2-Dibromo-3-Chloropropane | 10.0 | 8.17 | 82 | 50-132 | |
| 1,2-Dichlorobenzene | 10.0 | 9.02 | 90 | 79-120 | |
| 1,2-Dichloroethane | 10.0 | 9.32 | 93 | 80-120 | |
| 1,2-Dichloropropane | 10.0 | 9.80 | 98 | 78-124 | |
| 1,3-Dichlorobenzene | 10.0 | 9.06 | 91 | 79-120 | |
| 1,3-Dichloropropane | 10.0 | 9.36 | 94 | 80-120 | |
| 1,4-Dichlorobenzene | 10.0 | 9.07 | 91 | 79-120 | |
| 2,2-Dichloropropane | 10.0 | 7.92 | 79 | 71-122 | |
| 2-Butanone (MEK) | 20.0 | 21.5 | 108 | 56-138 | |
| 2-Chloroethyl vinyl ether | 10.0 | 9.80 J | 98 | 54-139 | |
| 2-Chlorotoluene | 10.0 | 9.14 | 91 | 80-120 | |
| 2-Hexanone | 20.0 | 18.5 | 92 | 55-141 | |
| 4-Chlorotoluene | 10.0 | 9.06 | 91 | 80-120 | |
| p-Isopropyltoluene | 10.0 | 9.23 | 92 | 77-120 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 19.6 | 98 | 64-135 | |
| Acetone | 20.0 | 19.1 | 95 | 34-148 | |
| Benzene | 10.0 | 9.39 | 94 | 80-120 | |
| Bromobenzene | 10.0 | 9.14 | 91 | 80-120 | |
| Bromoform | 10.0 | 6.29 | 63 | 56-122 | |
| Bromomethane | 10.0 | 6.43 | 64 | 38-132 | |
| Carbon disulfide | 10.0 | 9.43 | 94 | 65-144 | |
| Carbon tetrachloride | 10.0 | 7.89 | 79 | 77-131 | |
| Chlorobenzene | 10.0 | 9.32 | 93 | 80-120 | |
| Dibromochloromethane | 10.0 | 7.97 | 80 | 74-120 | |
| Chloroethane | 10.0 | 7.54 | 75 | 36-126 | |
| Chloroform | 10.0 | 9.57 | 96 | 80-120 | |
| Chloromethane | 10.0 | 9.55 | 96 | 48-133 | |
| cis-1,2-Dichloroethene | 10.0 | 9.75 | 98 | 79-120 | |
| cis-1,3-Dichloropropene | 10.0 | 9.84 | 98 | 74-126 | |
| Dibromomethane | 10.0 | 9.81 | 98 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC5593.D

Lab ID: LCS 240-212893/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 10.0 | 9.83 | 98 | 80-120 | |
| Dichlorodifluoromethane | 10.0 | 7.57 | 76 | 23-136 | |
| Bromodichloromethane | 10.0 | 8.91 | 89 | 80-120 | |
| tert-Butyl alcohol (TBA) | 100 | 79.5 | 79 | 38-144 | |
| Ethylbenzene | 10.0 | 9.45 | 95 | 80-120 | |
| Hexachlorobutadiene | 10.0 | 7.63 | 76 | 62-120 | |
| Isopropylbenzene | 10.0 | 9.47 | 95 | 77-120 | |
| Methyl tert-butyl ether | 10.0 | 9.50 | 95 | 69-121 | |
| Methylene Chloride | 10.0 | 8.56 | 86 | 77-129 | |
| m-Xylene & p-Xylene | 10.0 | 9.38 | 94 | 80-120 | |
| Naphthalene | 10.0 | 9.11 | 91 | 53-121 | |
| n-Butylbenzene | 10.0 | 8.99 | 90 | 74-120 | |
| N-Propylbenzene | 10.0 | 9.29 | 93 | 79-123 | |
| o-Xylene | 10.0 | 9.59 | 96 | 80-120 | |
| sec-Butylbenzene | 10.0 | 9.27 | 93 | 77-120 | |
| Styrene | 10.0 | 9.62 | 96 | 76-122 | |
| Tetrachloroethene | 10.0 | 9.39 | 94 | 78-121 | |
| Toluene | 10.0 | 9.29 | 93 | 80-120 | |
| trans-1,2-Dichloroethene | 10.0 | 9.74 | 97 | 80-124 | |
| trans-1,3-Dichloropropene | 10.0 | 9.12 | 91 | 75-131 | |
| Trichloroethene | 10.0 | 9.82 | 98 | 80-121 | |
| Trichlorofluoromethane | 10.0 | 9.18 | 92 | 61-133 | |
| Vinyl acetate | 10.0 | 8.16 | 82 | 27-160 | |
| Vinyl chloride | 10.0 | 9.13 | 91 | 52-121 | |
| Xylenes, Total | 20.0 | 19.0 | 95 | 80-120 | |
| tert-Butylbenzene | 10.0 | 9.36 | 94 | 77-121 | |
| 1,2-Dibromoethane | 10.0 | 9.65 | 96 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 147327.D

Lab ID: 240-59482-2 MS  Client ID: F-VB-UST-P-7 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 54.2 | 0.58 U | 44.5 | 82 | 34-135 | |
| 1,1,1-Trichloroethane | 54.2 | 0.76 U | 50.6 | 93 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 54.2 | 0.36 U | 43.6 | 80 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 54.2 | 0.49 U | 52.2 | 96 | 50-147 | |
| 1,1-Dichloroethane | 54.2 | 0.40 U | 45.9 | 85 | 54-122 | |
| 1,1-Dichloroethene | 54.2 | 0.93 U | 48.2 | 89 | 49-157 | |
| 1,1-Dichloropropene | 54.2 | 0.43 U | 50.4 | 93 | 50-122 | |
| 1,2,3-Trichlorobenzene | 54.2 | 0.58 U | 22.8 | 42 | 10-120 | |
| 1,2,3-Trichloropropane | 54.2 | 0.91 U | 48.1 | 89 | 32-174 | |
| 1,2,4-Trichlorobenzene | 54.2 | 0.44 U | 23.3 | 43 | 10-120 | |
| 1,2,4-Trimethylbenzene | 54.2 | 1.9 J | 37.0 | 65 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 54.2 | 3.2 U | 40.1 | 74 | 10-153 | |
| 1,2-Dichlorobenzene | 54.2 | 0.40 U | 30.9 | 57 | 17-122 | |
| 1,2-Dichloroethane | 54.2 | 0.54 U | 44.3 | 82 | 49-123 | |
| 1,2-Dichloropropane | 54.2 | 0.16 U | 47.0 | 87 | 61-120 | |
| 1,3-Dichlorobenzene | 54.2 | 0.80 U | 29.9 | 55 | 16-126 | |
| 1,3-Dichloropropane | 54.2 | 0.47 U | 41.4 | 76 | 54-128 | |
| 1,4-Dichlorobenzene | 54.2 | 0.33 U | 27.7 | 51 | 15-121 | |
| 2,2-Dichloropropane | 54.2 | 0.51 U | 46.0 | 85 | 49-132 | |
| 2-Butanone (MEK) | 108 | 4.8 J | 89.2 | 78 | 30-143 | |
| 2-Chloroethyl vinyl ether | 54.2 | 0.75 U | 33.0 J | 61 | 10-162 | |
| 2-Chlorotoluene | 54.2 | 0.35 U | 35.4 | 65 | 11-162 | |
| 2-Hexanone | 108 | 0.79 U | 90.1 | 83 | 37-147 | |
| 4-Chlorotoluene | 54.2 | 0.30 U | 32.1 | 59 | 15-149 | |
| p-Isopropyltoluene | 54.2 | 0.48 U | 34.2 | 63 | 10-165 | |
| 4-Methyl-2-pentanone (MIBK) | 108 | 1.4 U | 97.1 | 90 | 43-147 | |
| Acetone | 108 | 36 | 102 | 61 | 24-140 | |
| Benzene | 54.2 | 1.0 U | 45.9 | 85 | 53-120 | |
| Bromobenzene | 54.2 | 0.38 U | 35.4 | 65 | 24-144 | |
| Bromoform | 54.2 | 0.28 U | 42.3 | 78 | 18-129 | |
| Bromomethane | 21.7 | 0.49 U | 16.0 | 74 | 33-130 | |
| Carbon disulfide | 54.2 | 0.69 U | 43.4 | 80 | 20-151 | |
| Carbon tetrachloride | 54.2 | 0.76 U | 54.2 | 100 | 32-137 | |
| Chlorobenzene | 54.2 | 0.62 U | 38.3 | 71 | 37-120 | |
| Dibromochloromethane | 54.2 | 0.43 U | 42.4 | 78 | 29-135 | |
| Chloroethane | 21.7 | 0.51 U | 17.2 | 79 | 45-120 | |
| Chloroform | 54.2 | 0.42 U | 45.3 | 83 | 53-120 | |
| Chloromethane | 21.7 | 0.92 U | 18.5 | 85 | 34-120 | |
| cis-1,2-Dichloroethene | 54.2 | 0.33 U | 43.2 | 80 | 50-120 | |
| cis-1,3-Dichloropropene | 54.2 | 0.83 U | 42.5 | 78 | 27-133 | |
| Dibromomethane | 54.2 | 0.49 U | 43.3 | 80 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 147327.D

Lab ID: 240-59482-2 MS Client ID: F-VB-UST-P-7 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Bromochloromethane | 54.2 | 0.38 U | 41.7 | 77 | 53-120 | |
| Dichlorodifluoromethane | 21.7 | 0.33 U | 20.5 | 95 | 17-120 | |
| Bromodichloromethane | 54.2 | 0.34 U | 46.2 | 85 | 35-132 | |
| tert-Butyl alcohol (TBA) | 542 | 6.7 U | 454 | 84 | 60-125 | |
| Ethylbenzene | 54.2 | 0.66 J | 40.5 | 73 | 30-131 | |
| Hexachlorobutadiene | 54.2 | 0.47 U | 24.2 | 45 | 10-131 | |
| Isopropylbenzene | 54.2 | 0.24 J | 39.5 | 72 | 21-134 | |
| Methyl tert-butyl ether | 54.2 | 0.50 U | 46.4 | 85 | 51-157 | |
| Methylene Chloride | 54.2 | 0.86 U | 40.6 | 75 | 54-120 | |
| m-Xylene & p-Xylene | 54.2 | 2.9 J | 42.1 | 72 | 29-131 | |
| Naphthalene | 54.2 | 0.61 J | 31.1 | 56 | 10-124 | |
| n-Butylbenzene | 54.2 | 0.69 U | 29.5 | 54 | 10-148 | |
| N-Propylbenzene | 54.2 | 0.70 U | 36.8 | 68 | 10-178 | |
| o-Xylene | 54.2 | 1.2 J | 42.2 | 76 | 29-134 | |
| sec-Butylbenzene | 54.2 | 1.1 U | 35.7 | 66 | 10-172 | |
| Styrene | 54.2 | 0.47 U | 37.2 | 69 | 27-127 | |
| Tetrachloroethene | 54.2 | 0.93 U | 42.6 | 78 | 31-135 | |
| Toluene | 54.2 | 1.6 J | 45.1 | 80 | 39-129 | |
| trans-1,2-Dichloroethene | 54.2 | 0.37 U | 44.0 | 81 | 50-123 | |
| trans-1,3-Dichloropropene | 54.2 | 0.34 U | 37.3 | 69 | 28-137 | |
| Trichloroethene | 54.2 | 0.44 U | 44.7 | 82 | 10-177 | |
| Trichlorofluoromethane | 21.7 | 0.34 U | 20.8 | 96 | 36-142 | |
| Vinyl acetate | 54.2 | 0.77 U | 32.9 | 61 | 70-130 | F1 |
| Vinyl chloride | 21.7 | 0.35 U | 18.3 | 84 | 42-120 | |
| Xylenes, Total | 108 | 4.1 J | 84.3 | 74 | 30-131 | |
| tert-Butylbenzene | 54.2 | 0.78 U | 38.2 | 70 | 10-163 | |
| 1,2-Dibromoethane | 54.2 | 0.62 U | 42.6 | 78 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC5610.D

Lab ID: 240-59545-B-2 MS Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 5000 | 220 U | 4250 | 85 | 69-122 | |
| 1,1,2,2-Tetrachloroethane | 5000 | 110 U | 4970 | 99 | 61-130 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 5000 | 230 U | 4400 | 88 | 44-140 | |
| 1,1-Dichloroethane | 5000 | 5600 | 10600 | 100 | 73-124 | |
| 1,1-Dichloroethene | 5000 | 230 U | 4740 | 95 | 67-124 | |
| 1,2,4-Trichlorobenzene | 5000 | 160 U | 3990 | 80 | 48-120 | |
| 1,2-Dibromo-3-Chloropropane | 5000 | 410 U | 3890 | 78 | 42-130 | |
| 1,2-Dichlorobenzene | 5000 | 130 U | 4420 | 88 | 67-118 | |
| 1,2-Dichloroethane | 5000 | 120 J | 4970 | 99 | 74-125 | |
| 1,2-Dichloropropane | 5000 | 130 U | 4850 | 97 | 73-122 | |
| 1,3-Dichlorobenzene | 5000 | 95 U | 4440 | 89 | 65-120 | |
| 1,4-Dichlorobenzene | 5000 | 140 U | 4470 | 89 | 66-120 | |
| 2-Butanone (MEK) | 10000 | 660 J | 11300 | 107 | 49-132 | |
| 2-Hexanone | 10000 | 240 U | 10100 | 101 | 49-142 | |
| 4-Methyl-2-pentanone (MIBK) | 10000 | 500 U | 10400 | 104 | 58-136 | |
| Acetone | 10000 | 470 U | 8410 | 84 | 32-126 | |
| Benzene | 5000 | 180 U | 4720 | 94 | 73-121 | |
| Bromoform | 5000 | 280 U | 3020 | 60 | 45-121 | |
| Bromomethane | 5000 | 220 U | 2680 | 54 | 26-136 | |
| Carbon disulfide | 5000 | 190 U | 4490 | 90 | 54-144 | |
| Carbon tetrachloride | 5000 | 220 U | 3710 | 74 | 65-129 | |
| Chlorobenzene | 5000 | 130 U | 4670 | 93 | 72-120 | |
| Dibromochloromethane | 5000 | 220 U | 3910 | 78 | 62-122 | |
| Chloroethane | 5000 | 410 J | 3500 | 62 | 27-131 | |
| Chloroform | 5000 | 130 U | 4760 | 95 | 73-121 | |
| Chloromethane | 5000 | 220 U | 4750 | 95 | 39-134 | |
| cis-1,2-Dichloroethene | 5000 | 13000 | 17700 | 101 | 66-124 | |
| cis-1,3-Dichloropropene | 5000 | 230 U | 4550 | 91 | 60-120 | |
| Dichlorodifluoromethane | 5000 | 160 U | 3190 | 64 | 14-137 | |
| Bromodichloromethane | 5000 | 150 U | 4260 | 85 | 72-120 | |
| Ethylbenzene | 5000 | 130 U | 4710 | 94 | 68-121 | |
| Isopropylbenzene | 5000 | 180 U | 4710 | 94 | 61-122 | |
| Methyl tert-butyl ether | 5000 | 100 U | 4640 | 93 | 61-121 | |
| Methylene Chloride | 5000 | 170 U | 4360 | 87 | 70-124 | |
| m-Xylene & p-Xylene | 5000 | 120 U | 4700 | 94 | 66-123 | |
| o-Xylene | 5000 | 130 U | 4760 | 95 | 68-121 | |
| Styrene | 5000 | 230 U | 4790 | 96 | 64-126 | |
| Tetrachloroethene | 5000 | 160 U | 4540 | 91 | 59-125 | |
| Toluene | 5000 | 120 U | 4690 | 94 | 72-122 | |
| trans-1,2-Dichloroethene | 5000 | 500 | 5390 | 98 | 72-125 | |
| trans-1,3-Dichloropropene | 5000 | 280 U | 4330 | 87 | 58-132 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC5610.D

Lab ID: 240-59545-B-2 MS Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Trichloroethene | 5000 | 110 U | 4800 | 96 | 61-129 | |
| Trichlorofluoromethane | 5000 | 250 U | 3960 | 79 | 49-133 | |
| Vinyl chloride | 5000 | 4200 | 8920 | 95 | 44-122 | |
| Xylenes, Total | 10000 | 260 U | 9460 | 95 | 67-122 | |
| 1,2-Dibromoethane | 5000 | 160 U | 5040 | 101 | 69-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 147328.D

Lab ID: 240-59482-2 MSD Client ID: F-VB-UST-P-7 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 56.0 | 41.7 | 74 | 7 | 30 | 34-135 | |
| 1,1,1-Trichloroethane | 56.0 | 55.0 | 98 | 8 | 30 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 56.0 | 47.6 | 85 | 9 | 30 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 56.0 | 58.3 | 104 | 11 | 30 | 50-147 | |
| 1,1-Dichloroethane | 56.0 | 50.9 | 91 | 10 | 30 | 54-122 | |
| 1,1-Dichloroethene | 56.0 | 56.9 | 102 | 17 | 30 | 49-157 | |
| 1,1-Dichloropropene | 56.0 | 51.9 | 93 | 3 | 30 | 50-122 | |
| 1,2,3-Trichlorobenzene | 56.0 | 19.0 | 34 | 18 | 30 | 10-120 | |
| 1,2,3-Trichloropropane | 56.0 | 50.5 | 90 | 5 | 30 | 32-174 | |
| 1,2,4-Trichlorobenzene | 56.0 | 17.4 | 31 | 29 | 30 | 10-120 | |
| 1,2,4-Trimethylbenzene | 56.0 | 27.3 | 45 | 30 | 30 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 56.0 | 48.9 | 87 | 20 | 30 | 10-153 | |
| 1,2-Dichlorobenzene | 56.0 | 23.5 | 42 | 27 | 30 | 17-122 | |
| 1,2-Dichloroethane | 56.0 | 49.6 | 89 | 11 | 30 | 49-123 | |
| 1,2-Dichloropropane | 56.0 | 48.5 | 87 | 3 | 30 | 61-120 | |
| 1,3-Dichlorobenzene | 56.0 | 20.8 | 37 | 36 | 30 | 16-126 | F2 |
| 1,3-Dichloropropane | 56.0 | 43.5 | 78 | 5 | 30 | 54-128 | |
| 1,4-Dichlorobenzene | 56.0 | 20.0 | 36 | 32 | 30 | 15-121 | F2 |
| 2,2-Dichloropropane | 56.0 | 50.4 | 90 | 9 | 30 | 49-132 | |
| 2-Butanone (MEK) | 112 | 106 | 91 | 18 | 30 | 30-143 | |
| 2-Chloroethyl vinyl ether | 56.0 | 39.9 J | 71 | 19 | 30 | 10-162 | |
| 2-Chlorotoluene | 56.0 | 24.1 | 43 | 38 | 30 | 11-162 | F2 |
| 2-Hexanone | 112 | 103 | 92 | 13 | 30 | 37-147 | |
| 4-Chlorotoluene | 56.0 | 21.3 | 38 | 40 | 30 | 15-149 | F2 |
| p-Isopropyltoluene | 56.0 | 20.0 | 36 | 52 | 30 | 10-165 | F2 |
| 4-Methyl-2-pentanone (MIBK) | 112 | 117 | 104 | 19 | 30 | 43-147 | |
| Acetone | 112 | 125 | 80 | 20 | 30 | 24-140 | |
| Benzene | 56.0 | 47.6 | 85 | 4 | 30 | 53-120 | |
| Bromobenzene | 56.0 | 28.4 | 51 | 22 | 30 | 24-144 | |
| Bromoform | 56.0 | 42.8 | 76 | 1 | 30 | 18-129 | |
| Bromomethane | 22.4 | 17.9 | 80 | 11 | 30 | 33-130 | |
| Carbon disulfide | 56.0 | 45.7 | 82 | 5 | 30 | 20-151 | |
| Carbon tetrachloride | 56.0 | 56.7 | 101 | 5 | 30 | 32-137 | |
| Chlorobenzene | 56.0 | 31.6 | 56 | 19 | 30 | 37-120 | |
| Dibromochloromethane | 56.0 | 43.6 | 78 | 3 | 30 | 29-135 | |
| Chloroethane | 22.4 | 19.1 | 85 | 10 | 30 | 45-120 | |
| Chloroform | 56.0 | 49.0 | 87 | 8 | 30 | 53-120 | |
| Chloromethane | 22.4 | 22.6 | 101 | 20 | 30 | 34-120 | |
| cis-1,2-Dichloroethene | 56.0 | 47.7 | 85 | 10 | 30 | 50-120 | |
| cis-1,3-Dichloropropene | 56.0 | 41.8 | 75 | 2 | 30 | 27-133 | |
| Dibromomethane | 56.0 | 46.2 | 83 | 7 | 30 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 147328.D

Lab ID: 240-59482-2 MSD Client ID: F-VB-UST-P-7 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Bromochloromethane | 56.0 | 46.8 | 84 | 12 | 30 | 53-120 | |
| Dichlorodifluoromethane | 22.4 | 24.1 | 108 | 16 | 30 | 17-120 | |
| Bromodichloromethane | 56.0 | 47.1 | 84 | 2 | 30 | 35-132 | |
| tert-Butyl alcohol (TBA) | 560 | 541 | 97 | 17 | 30 | 60-125 | |
| Ethylbenzene | 56.0 | 31.4 | 55 | 25 | 30 | 30-131 | |
| Hexachlorobutadiene | 56.0 | 12.2 | 22 | 66 | 30 | 10-131 | F2 |
| Isopropylbenzene | 56.0 | 28.6 | 51 | 32 | 30 | 21-134 | F2 |
| Methyl tert-butyl ether | 56.0 | 53.0 | 95 | 13 | 30 | 51-157 | |
| Methylene Chloride | 56.0 | 45.3 | 81 | 11 | 30 | 54-120 | |
| m-Xylene & p-Xylene | 56.0 | 34.9 | 57 | 19 | 30 | 29-131 | |
| Naphthalene | 56.0 | 31.3 | 55 | 1 | 30 | 10-124 | |
| n-Butylbenzene | 56.0 | 15.6 | 28 | 62 | 30 | 10-148 | F2 |
| N-Propylbenzene | 56.0 | 23.4 | 42 | 45 | 30 | 10-178 | F2 |
| o-Xylene | 56.0 | 33.9 | 58 | 22 | 30 | 29-134 | |
| sec-Butylbenzene | 56.0 | 21.4 | 38 | 50 | 30 | 10-172 | F2 |
| Styrene | 56.0 | 28.1 | 50 | 28 | 30 | 27-127 | |
| Tetrachloroethene | 56.0 | 34.7 | 62 | 20 | 30 | 31-135 | |
| Toluene | 56.0 | 40.2 | 69 | 11 | 30 | 39-129 | |
| trans-1,2-Dichloroethene | 56.0 | 48.3 | 86 | 9 | 30 | 50-123 | |
| trans-1,3-Dichloropropene | 56.0 | 38.8 | 69 | 4 | 30 | 28-137 | |
| Trichloroethene | 56.0 | 44.3 | 79 | 1 | 30 | 10-177 | |
| Trichlorofluoromethane | 22.4 | 24.7 | 110 | 17 | 30 | 36-142 | |
| Vinyl acetate | 56.0 | 43.6 | 78 | 28 | 30 | 70-130 | |
| Vinyl chloride | 22.4 | 21.8 | 97 | 18 | 30 | 42-120 | |
| Xylenes, Total | 112 | 68.8 | 58 | 20 | 30 | 30-131 | |
| tert-Butylbenzene | 56.0 | 25.3 | 45 | 40 | 30 | 10-163 | F2 |
| 1,2-Dibromoethane | 56.0 | 43.9 | 78 | 3 | 30 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC5611.D

Lab ID: 240-59545-B-2 MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 5000 | 4480 | 90 | 5 | 14 | 69-122 | |
| 1,1,2,2-Tetrachloroethane | 5000 | 4970 | 99 | 0 | 18 | 61-130 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 5000 | 5180 | 104 | 16 | 35 | 44-140 | |
| 1,1-Dichloroethane | 5000 | 10800 | 106 | 3 | 14 | 73-124 | |
| 1,1-Dichloroethene | 5000 | 5020 | 100 | 6 | 24 | 67-124 | |
| 1,2,4-Trichlorobenzene | 5000 | 4280 | 86 | 7 | 28 | 48-120 | |
| 1,2-Dibromo-3-Chloropropane | 5000 | 3860 | 77 | 1 | 24 | 42-130 | |
| 1,2-Dichlorobenzene | 5000 | 4610 | 92 | 4 | 15 | 67-118 | |
| 1,2-Dichloroethane | 5000 | 4960 | 99 | 0 | 24 | 74-125 | |
| 1,2-Dichloropropane | 5000 | 4910 | 98 | 1 | 15 | 73-122 | |
| 1,3-Dichlorobenzene | 5000 | 4550 | 91 | 2 | 15 | 65-120 | |
| 1,4-Dichlorobenzene | 5000 | 4580 | 92 | 2 | 16 | 66-120 | |
| 2-Butanone (MEK) | 10000 | 9840 | 92 | 14 | 19 | 49-132 | |
| 2-Hexanone | 10000 | 8950 | 89 | 12 | 27 | 49-142 | |
| 4-Methyl-2-pentanone (MIBK) | 10000 | 9860 | 99 | 5 | 32 | 58-136 | |
| Acetone | 10000 | 8100 | 81 | 4 | 28 | 32-126 | |
| Benzene | 5000 | 4820 | 96 | 2 | 13 | 73-121 | |
| Bromoform | 5000 | 3200 | 64 | 6 | 19 | 45-121 | |
| Bromomethane | 5000 | 5100 | 102 | 62 | 35 | 26-136 | F2 |
| Carbon disulfide | 5000 | 4730 | 95 | 5 | 34 | 54-144 | |
| Carbon tetrachloride | 5000 | 4200 | 84 | 12 | 20 | 65-129 | |
| Chlorobenzene | 5000 | 4690 | 94 | 0 | 15 | 72-120 | |
| Dibromochloromethane | 5000 | 3960 | 79 | 1 | 19 | 62-122 | |
| Chloroethane | 5000 | 5260 | 97 | 40 | 35 | 27-131 | F2 |
| Chloroform | 5000 | 4950 | 99 | 4 | 17 | 73-121 | |
| Chloromethane | 5000 | 4860 | 97 | 2 | 20 | 39-134 | |
| cis-1,2-Dichloroethene | 5000 | 18400 | 116 | 4 | 22 | 66-124 | |
| cis-1,3-Dichloropropene | 5000 | 4610 | 92 | 1 | 21 | 60-120 | |
| Dichlorodifluoromethane | 5000 | 4070 | 81 | 24 | 34 | 14-137 | |
| Bromodichloromethane | 5000 | 4410 | 88 | 3 | 19 | 72-120 | |
| Ethylbenzene | 5000 | 4850 | 97 | 3 | 16 | 68-121 | |
| Isopropylbenzene | 5000 | 4950 | 99 | 5 | 20 | 61-122 | |
| Methyl tert-butyl ether | 5000 | 4740 | 95 | 2 | 12 | 61-121 | |
| Methylene Chloride | 5000 | 4590 | 92 | 5 | 14 | 70-124 | |
| m-Xylene & p-Xylene | 5000 | 4800 | 96 | 2 | 15 | 66-123 | |
| o-Xylene | 5000 | 4820 | 96 | 1 | 14 | 68-121 | |
| Styrene | 5000 | 4900 | 98 | 2 | 15 | 64-126 | |
| Tetrachloroethene | 5000 | 4620 | 92 | 2 | 20 | 59-125 | |
| Toluene | 5000 | 4620 | 92 | 2 | 15 | 72-122 | |
| trans-1,2-Dichloroethene | 5000 | 5740 | 105 | 6 | 25 | 72-125 | |
| trans-1,3-Dichloropropene | 5000 | 4270 | 85 | 1 | 22 | 58-132 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC5611.D

Lab ID: 240-59545-B-2 MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Trichloroethene | 5000 | 4960 | 99 | 3 | 14 | 61-129 | |
| Trichlorofluoromethane | 5000 | 4740 | 95 | 18 | 25 | 49-133 | |
| Vinyl chloride | 5000 | 9380 | 104 | 5 | 35 | 44-122 | |
| Xylenes, Total | 10000 | 9620 | 96 | 2 | 14 | 67-122 | |
| 1,2-Dibromoethane | 5000 | 4900 | 98 | 3 | 24 | 69-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab File ID: 147321.D Lab Sample ID: MB 240-212606/24

Matrix: Solid Heated Purge:(Y/N) Y

Instrument ID: A3UX14 Date Analyzed: 12/30/2015 20:00

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-212606/23 | 147320.D | 12/30/2015 19:39 |
| | LCS 240-212606/25 | 147322.D | 12/30/2015 20:22 |
| F-VB-UST-P-7 | 240-59482-2 | 147326.D | 12/30/2015 21:48 |
| F-VB-UST-P-7 MS | 240-59482-2 MS | 147327.D | 12/30/2015 22:09 |
| F-VB-UST-P-7 MSD | 240-59482-2 MSD | 147328.D | 12/30/2015 22:31 |
| F-VB-UST-P-8 | 240-59482-3 | 147330.D | 12/30/2015 23:14 |

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab File ID: UXC5596.D Lab Sample ID: MB 240-212893/8

Matrix: Water Heated Purge:(Y/N) N

Instrument ID: A3UX15 Date Analyzed: 01/04/2016 11:26

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-212893/5 | UXC5593.D | 01/04/2016 10:18 |
| TB-122815 | 240-59482-1 | UXC5597.D | 01/04/2016 12:07 |
| | 240-59545-B-2 MS | UXC5610.D | 01/04/2016 17:00 |
| | 240-59545-B-2 MSD | UXC5611.D | 01/04/2016 17:23 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab File ID: BFB Injection Date:

Instrument ID: BFB Injection Time:

Lab File ID: DFTPP Injection Date:

Instrument ID: DFTPP Injection Time:

Analysis Batch No.:

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|---|---|---|
| | | |

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | IC 240-212606/3 | 147300.D | 12/30/2015 | 12:29 |
| | IC 240-212606/4 | 147301.D | 12/30/2015 | 12:50 |
| | IC 240-212606/5 | 147302.D | 12/30/2015 | 13:12 |
| | ICIS 240-212606/6 | 147303.D | 12/30/2015 | 13:33 |
| | IC 240-212606/7 | 147304.D | 12/30/2015 | 13:55 |
| | IC 240-212606/8 | 147305.D | 12/30/2015 | 14:16 |
| | IC 240-212606/9 | 147306.D | 12/30/2015 | 14:38 |
| | IC 240-212606/10 | 147307.D | 12/30/2015 | 14:59 |
| | IC 240-212606/11 | 147308.D | 12/30/2015 | 15:21 |
| | IC 240-212606/12 | 147309.D | 12/30/2015 | 15:42 |
| | IC 240-212606/13 | 147310.D | 12/30/2015 | 16:04 |
| | IC 240-212606/14 | 147311.D | 12/30/2015 | 16:25 |
| | IC 240-212606/15 | 147312.D | 12/30/2015 | 16:47 |
| | IC 240-212606/16 | 147313.D | 12/30/2015 | 17:08 |
| | IC 240-212606/17 | 147314.D | 12/30/2015 | 17:30 |
| | IC 240-212606/18 | 147315.D | 12/30/2015 | 17:51 |
| | IC 240-212606/19 | 147316.D | 12/30/2015 | 18:13 |
| | IC 240-212606/20 | 147317.D | 12/30/2015 | 18:34 |
| | ICV 240-212606/21 | 147318.D | 12/30/2015 | 18:56 |
| | ICV 240-212606/22 | 147319.D | 12/30/2015 | 19:17 |
| | LCS 240-212606/23 | 147320.D | 12/30/2015 | 19:39 |
| | MB 240-212606/24 | 147321.D | 12/30/2015 | 20:00 |
| | LCS 240-212606/25 | 147322.D | 12/30/2015 | 20:22 |
| F-VB-UST-P-7 | 240-59482-2 | 147326.D | 12/30/2015 | 21:48 |
| F-VB-UST-P-7 MS | 240-59482-2 MS | 147327.D | 12/30/2015 | 22:09 |
| F-VB-UST-P-7 MSD | 240-59482-2 MSD | 147328.D | 12/30/2015 | 22:31 |
| F-VB-UST-P-8 | 240-59482-3 | 147330.D | 12/30/2015 | 23:14 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab File ID: BFB496.D BFB Injection Date: 12/14/2015

Instrument ID: A3UX15 BFB Injection Time: 12:30

Analysis Batch No.: 210738

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 16.9 | |
| 75 | 30.0 - 60.0 % of mass 95 | 48.5 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.6 | |
| 173 | Less than 2.0 % of mass 174 | 0.4 | (0.6) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 81.1 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.4 | (6.7) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 78.6 | (96.9) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 5.2 | (6.6) 2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD8260 240-210738/2 | UXC5222A.D | 12/14/2015 | 13:08 |
| | STD8260 240-210738/3 | UXC5223.D | 12/14/2015 | 13:31 |
| | STD8260 240-210738/4 | UXC5224.D | 12/14/2015 | 13:54 |
| | STD8260 240-210738/5 | UXC5225.D | 12/14/2015 | 14:16 |
| | STD8260 240-210738/6 | UXC5226.D | 12/14/2015 | 14:39 |
| | STD8260 240-210738/7 | UXC5227.D | 12/14/2015 | 15:01 |
| | ICV 240-210738/14 | UXC5228.D | 12/14/2015 | 15:24 |
| | STD 240-210738/8 | UXC5229.D | 12/14/2015 | 15:47 |
| | STD 240-210738/9 | UXC5230.D | 12/14/2015 | 16:09 |
| | STD 240-210738/10 | UXC5231.D | 12/14/2015 | 16:32 |
| | STD 240-210738/11 | UXC5232.D | 12/14/2015 | 16:55 |
| | STD 240-210738/12 | UXC5233.D | 12/14/2015 | 17:17 |
| | STD 240-210738/13 | UXC5234.D | 12/14/2015 | 17:40 |
| | ICV 240-210738/15 | UXC5235.D | 12/14/2015 | 18:03 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab File ID: BFB505.D  BFB Injection Date: 01/04/2016

Instrument ID: A3UX15  BFB Injection Time: 09:30

Analysis Batch No.: 212893

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 16.9 | |
| 75 | 30.0 - 60.0 % of mass 95 | 48.8 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.6 | |
| 173 | Less than 2.0 % of mass 174 | 0.4 | (0.5) 1 |
| 174 | 50.0 - 120.00 % of mass 95 | 77.1 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.2 | (6.7) 1 |
| 176 | 95.0 - 101.0 % of mass 174 | 73.6 | (95.4) 1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.7 | (6.4) 2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-212893/2 | UXC5592.D | 01/04/2016 | 09:55 |
| | LCS 240-212893/5 | UXC5593.D | 01/04/2016 | 10:18 |
| | CCV 240-212893/3 | UXC5594.D | 01/04/2016 | 10:41 |
| | MB 240-212893/8 | UXC5596.D | 01/04/2016 | 11:26 |
| TB-122815 | 240-59482-1 | UXC5597.D | 01/04/2016 | 12:07 |
| | 240-59545-B-2 MS | UXC5610.D | 01/04/2016 | 17:00 |
| | 240-59545-B-2 MSD | UXC5611.D | 01/04/2016 | 17:23 |

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Sample No.: ICIS 240-212606/6  Date Analyzed: 12/30/2015 13:33

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 147303.D  Heated Purge: (Y/N) Y

Calibration ID: 32237

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 1280920 | 6.64 | 787274 | 9.36 | 374460 | 11.34 |
| UPPER LIMIT | | 2561840 | 7.14 | 1574548 | 9.86 | 748920 | 11.84 |
| LOWER LIMIT | | 640460 | 6.14 | 393637 | 8.86 | 187230 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-212606/21 | | 1152357 | 6.64 | 718027 | 9.36 | 346347 | 11.34 |
| ICV 240-212606/22 | | 1170616 | 6.64 | 666787 | 9.36 | 354139 | 11.34 |
| LCS 240-212606/23 | | 1187069 | 6.64 | 689945 | 9.36 | 347278 | 11.34 |
| MB 240-212606/24 | | 1153788 | 6.64 | 711695 | 9.36 | 363540 | 11.34 |
| LCS 240-212606/25 | | 1186477 | 6.64 | 716186 | 9.36 | 360250 | 11.34 |
| 240-59482-2 | F-VB-UST-P-7 | 1080515 | 6.64 | 664224 | 9.36 | 345118 | 11.34 |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 1170351 | 6.64 | 730231 | 9.36 | 366151 | 11.34 |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 1015563 | 6.64 | 632992 | 9.36 | 315004 | 11.34 |
| 240-59482-3 | F-VB-UST-P-8 | 1057314 | 6.64 | 632789 | 9.36 | 325058 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Sample No.: STD8260 240-210738/4  Date Analyzed: 12/14/2015 13:54

Instrument ID: A3UX15  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UXC5224.D  Heated Purge: (Y/N) N

Calibration ID: 32097

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 1817560 | 4.97 | 1323725 | 7.64 | 680775 | 9.85 |
| UPPER LIMIT | | 3635120 | 5.47 | 2647450 | 8.14 | 1361550 | 10.35 |
| LOWER LIMIT | | 908780 | 4.47 | 661863 | 7.14 | 340388 | 9.35 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-210738/14 | | 1720163 | 4.98 | 1275243 | 7.64 | 670998 | 9.85 |
| ICV 240-210738/15 | | 1705447 | 4.97 | 1245348 | 7.64 | 646436 | 9.85 |
| CCVIS 240-212893/2 | | 1754912 | 4.98 | 1366553 | 7.64 | 728534 | 9.85 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Sample No.: CCVIS 240-212893/2  Date Analyzed: 01/04/2016 09:55

Instrument ID: A3UX15  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UXC5592.D  Heated Purge: (Y/N) N

Calibration ID: 32101

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1754912 | 4.98 | 1366553 | 7.64 | 728534 | 9.85 |
| UPPER LIMIT | | 3509824 | 5.48 | 2733106 | 8.14 | 1457068 | 10.35 |
| LOWER LIMIT | | 877456 | 4.48 | 683277 | 7.14 | 364267 | 9.35 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| LCS 240-212893/5 | | 1698190 | 4.97 | 1325194 | 7.64 | 719950 | 9.85 |
| CCV 240-212893/3 | | 1694967 | 4.97 | 1320936 | 7.64 | 698544 | 9.85 |
| MB 240-212893/8 | | 1703116 | 4.98 | 1282963 | 7.64 | 684171 | 9.85 |
| 240-59482-1 | TB-122815 | 1675391 | 4.97 | 1275724 | 7.64 | 666155 | 9.85 |
| 240-59545-B-2 MS | | 1682893 | 4.97 | 1290316 | 7.64 | 702033 | 9.85 |
| 240-59545-B-2 MSD | | 1618113 | 4.98 | 1294288 | 7.64 | 708761 | 9.85 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: TB-122815 | Lab Sample ID: 240-59482-1

Matrix: Water | Lab File ID: UXC5597.D

Analysis Method: 8260C | Date Collected: 12/28/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 12:07

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.28 | U | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.44 | U | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.22 | U | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.30 | U | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 0.45 | U | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.37 | U | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.44 | U | 1.0 | 0.44 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.47 | U | 5.0 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.32 | U | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.41 | U | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.82 | U | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.25 | U | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 0.23 | U | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 0.25 | U | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.19 | U | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 0.19 | U | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.27 | U | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 0.26 | U | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 1.5 | J B | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.59 | U | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 0.40 | U | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 0.48 | U | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 10 | 0.99 |
| 67-64-1 | Acetone | 1.8 | J | 10 | 0.94 |
| 71-43-2 | Benzene | 0.35 | U | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 0.35 | U | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 0.56 | U | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 0.44 | U | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.38 | U | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 0.43 | U | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 0.25 | U | 1.0 | 0.25 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: TB-122815 | Lab Sample ID: 240-59482-1

Matrix: Water | Lab File ID: UXC5597.D

Analysis Method: 8260C | Date Collected: 12/28/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 12:07

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 1.0 | 0.43 |
| 75-00-3 | Chloroethane | 0.32 | U | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 0.25 | U | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 0.44 | U | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.26 | U | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.46 | U | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 0.42 | U | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.50 | U | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.1 | J | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 0.25 | U | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 0.35 | U | 1.0 | 0.35 |
| 108-20-3 | Diisopropyl ether | 0.50 | U | 10 | 0.50 |
| 98-82-8 | Isopropylbenzene | 0.35 | U | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 0.20 | U | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 0.33 | U | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.24 | U | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 0.45 | U | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 0.31 | U | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 0.40 | U | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 0.25 | U | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 0.48 | U | 1.0 | 0.48 |
| 100-42-5 | Styrene | 0.45 | U | 1.0 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.30 | U | 5.0 | 0.30 |
| 637-92-3 | Ethyl t-butyl ether | 0.23 | U | 5.0 | 0.23 |
| 127-18-4 | Tetrachloroethene | 0.31 | U | 1.0 | 0.31 |
| 108-88-3 | Toluene | 0.23 | U | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.30 | U | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.56 | U | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 0.22 | U | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 0.49 | U | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 0.41 | U | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 0.29 | U | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 0.52 | U | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 0.41 | U | 1.0 | 0.41 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: TB-122815 | Lab Sample ID: 240-59482-1

Matrix: Water | Lab File ID: UXC5597.D

Analysis Method: 8260C | Date Collected: 12/28/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 12:07

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.32 | U | 1.0 | 0.32 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 102 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 93 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 112 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 99 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 | Lab Sample ID: 240-59482-2

Matrix: Solid | Lab File ID: 147326.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.26(g) | Date Analyzed: 12/30/2015 21:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.76 | U | 5.8 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 5.8 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.91 | U | 5.8 | 0.91 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.86 | J | 5.8 | 0.54 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.9 | J | 5.8 | 0.68 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 5.8 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.54 | U | 5.8 | 0.54 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U F2 | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.47 | U | 5.8 | 0.47 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U F2 | 5.8 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.8 | J | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.75 | U | 58 | 0.75 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U F2 | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U F2 | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.48 | U F2 | 5.8 | 0.48 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 36 | | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.69 | U | 5.8 | 0.69 |
| 56-23-5 | Carbon tetrachloride | 0.76 | U | 5.8 | 0.76 |
| 108-90-7 | Chlorobenzene | 0.62 | U | 5.8 | 0.62 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 | Lab Sample ID: 240-59482-2

Matrix: Solid | Lab File ID: 147326.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.26(g) | Date Analyzed: 12/30/2015 21:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 5.8 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.83 | U | 5.8 | 0.83 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 5.8 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.7 | U | 230 | 6.7 |
| 100-41-4 | Ethylbenzene | 0.66 | J | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.47 | U F2 | 5.8 | 0.47 |
| 108-20-3 | Diisopropyl ether | 0.69 | U | 12 | 0.69 |
| 98-82-8 | Isopropylbenzene | 0.24 | J F2 | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 2.9 | J | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.61 | J B | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 0.69 | U F2 | 5.8 | 0.69 |
| 103-65-1 | N-Propylbenzene | 0.70 | U F2 | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 1.2 | J | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U F2 | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.47 | U | 5.8 | 0.47 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.48 | U | 5.8 | 0.48 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 1.6 | J | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.77 | U F1 | 12 | 0.77 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 4.1 | J | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 0.78 | U F2 | 5.8 | 0.78 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 | Lab Sample ID: 240-59482-2

Matrix: Solid | Lab File ID: 147326.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.26(g) | Date Analyzed: 12/30/2015 21:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.62 | U | 5.8 | 0.62 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 91 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 93 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 88 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 100 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-8 | Lab Sample ID: 240-59482-3

Matrix: Solid | Lab File ID: 147330.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:30

Sample wt/vol: 5.43(g) | Date Analyzed: 12/30/2015 23:14

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.1 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.71 | U | 5.5 | 0.71 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.37 | U | 5.5 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 0.88 | U | 5.5 | 0.88 |
| 563-58-6 | 1,1-Dichloropropene | 0.41 | U | 5.5 | 0.41 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.86 | U | 5.5 | 0.86 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.50 | U | 5.5 | 0.50 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.42 | U | 5.5 | 0.42 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.64 | U | 5.5 | 0.64 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.37 | U | 5.5 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 0.50 | U | 5.5 | 0.50 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.76 | U | 5.5 | 0.76 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.48 | U | 5.5 | 0.48 |
| 78-93-3 | 2-Butanone (MEK) | 1.9 | J | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.70 | U | 55 | 0.70 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.75 | U | 22 | 0.75 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.5 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.57 | J | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 22 | 1.3 |
| 67-64-1 | Acetone | 9.4 | J | 22 | 6.3 |
| 71-43-2 | Benzene | 0.99 | U | 5.5 | 0.99 |
| 108-86-1 | Bromobenzene | 0.36 | U | 5.5 | 0.36 |
| 75-25-2 | Bromoform | 0.26 | U | 5.5 | 0.26 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.65 | U | 5.5 | 0.65 |
| 56-23-5 | Carbon tetrachloride | 0.71 | U | 5.5 | 0.71 |
| 108-90-7 | Chlorobenzene | 0.58 | U | 5.5 | 0.58 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-8 | Lab Sample ID: 240-59482-3

Matrix: Solid | Lab File ID: 147330.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:30

Sample wt/vol: 5.43(g) | Date Analyzed: 12/30/2015 23:14

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.1 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.41 | U | 5.5 | 0.41 |
| 75-00-3 | Chloroethane | 0.48 | U | 5.5 | 0.48 |
| 67-66-3 | Chloroform | 0.40 | U | 5.5 | 0.40 |
| 74-87-3 | Chloromethane | 0.87 | U | 5.5 | 0.87 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.78 | U | 5.5 | 0.78 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.36 | U | 5.5 | 0.36 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.3 | U | 220 | 6.3 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.65 | U | 11 | 0.65 |
| 98-82-8 | Isopropylbenzene | 2.0 | J | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.47 | U | 5.5 | 0.47 |
| 75-09-2 | Methylene Chloride | 0.81 | U | 5.5 | 0.81 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.88 | J B | 5.5 | 0.38 |
| 104-51-8 | n-Butylbenzene | 0.83 | J | 5.5 | 0.65 |
| 103-65-1 | N-Propylbenzene | 1.9 | J | 5.5 | 0.66 |
| 95-47-6 | o-Xylene | 0.63 | U | 5.5 | 0.63 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.5 | 1.0 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.47 | U | 5.5 | 0.47 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.88 | U | 5.5 | 0.88 |
| 108-88-3 | Toluene | 0.30 | U | 5.5 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.42 | U | 5.5 | 0.42 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.72 | U | 11 | 0.72 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.59 | U | 11 | 0.59 |
| 98-06-6 | tert-Butylbenzene | 0.74 | U | 5.5 | 0.74 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-8 | Lab Sample ID: 240-59482-3

Matrix: Solid | Lab File ID: 147330.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:30

Sample wt/vol: 5.43(g) | Date Analyzed: 12/30/2015 23:14

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.1 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.58 | U | 5.5 | 0.58 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 91 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 90 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 97 | | 58-123 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-212606/11 | 147308.D |
| Level 2 | IC 240-212606/10 | 147307.D |
| Level 3 | IC 240-212606/9 | 147306.D |
| Level 4 | IC 240-212606/8 | 147305.D |
| Level 5 | IC 240-212606/7 | 147304.D |
| Level 6 | ICIS 240-212606/6 | 147303.D |
| Level 7 | IC 240-212606/5 | 147302.D |
| Level 8 | IC 240-212606/4 | 147301.D |
| Level 9 | IC 240-212606/3 | 147300.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | 0.2522<br>0.3274 | 0.2622<br>0.3199 | 0.3084<br>0.3210 | 0.3205<br>0.3150 | 0.3246 | Ave | | 0.3057 | | | 0.1000 | 9.2 | | 15.0 | | | |
| Chloromethane | 0.4885<br>0.4863 | 0.4230<br>0.4845 | 0.4548<br>0.4778 | 0.4731<br>0.4486 | 0.4779 | Ave | | 0.4683 | | | 0.1000 | 4.7 | | 15.0 | | | |
| Vinyl chloride | 0.3094<br>0.3186 | 0.2798<br>0.3247 | 0.3148<br>0.3169 | 0.3087<br>0.3085 | 0.3161 | Ave | | 0.3108 | | | 0.1000 | 4.1 | | 15.0 | | | |
| Butadiene | 0.2974<br>0.3870 | 0.2971<br>0.3883 | 0.3583<br>0.3847 | 0.3906<br>0.3754 | 0.3828 | Ave | | 0.3624 | | | | 10.5 | | 15.0 | | | |
| Bromomethane | 0.1596<br>0.1219 | 0.1439<br>+++++ | 0.1309<br>+++++ | 0.1378<br>+++++ | 0.1335 | Ave | | 0.1379 | | | 0.0500 | 9.3 | | 15.0 | | | |
| Chloroethane | 0.2086<br>0.1635 | 0.1442<br>0.1527 | 0.1684<br>+++++ | 0.1696<br>+++++ | 0.1655 | Ave | | 0.1675 | | | | 12.1 | | 15.0 | | | |
| Dichlorofluoromethane | 0.3729<br>0.3474 | 0.3173<br>0.3305 | 0.3192<br>0.3405 | 0.3612<br>0.3272 | 0.3278 | Ave | | 0.3382 | | | | 5.6 | | 15.0 | | | |
| Trichlorofluoromethane | 0.3359<br>0.3669 | 0.2822<br>0.3627 | 0.3215<br>0.3573 | 0.3554<br>0.3540 | 0.3615 | Ave | | 0.3441 | | | 0.1000 | 7.9 | | 15.0 | | | |
| Ethyl ether | +++++<br>0.2212 | +++++<br>0.2173 | 0.2327<br>0.2086 | 0.2173<br>0.2046 | 0.2312 | Ave | | 0.2190 | | | | 4.8 | | 15.0 | | | |
| Acrolein | 0.0472<br>0.0341 | 0.0373<br>0.0351 | 0.0552<br>0.0318 | 0.0358<br>0.0321 | 0.0373 | Lin1 | 0.1371 | 0.0327 | | | | | | | 0.9960 | | 0.9900 |
| 1,1-Dichloroethene | 0.1970<br>0.2523 | 0.2302<br>0.2649 | 0.2739<br>0.2499 | 0.2592<br>0.2386 | 0.2682 | Ave | | 0.2482 | | | 0.1000 | 9.6 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.1681<br>0.2162 | 0.1916<br>0.2198 | 0.2273<br>0.2105 | 0.2084<br>0.2052 | 0.2286 | Ave | | 0.2084 | | | 0.0500 | 9.1 | | 15.0 | | | |
| Acetone | +++++<br>0.0865 | +++++<br>0.0863 | 0.1402<br>0.0797 | 0.1072<br>0.0753 | 0.1044 | Qua | 0.3866 | 0.0903 | -0.000031 | | 0.0500 | | | | 1.0000 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Iodomethane | 0.3774<br>0.3813 | 0.3660<br>0.3711 | 0.3712<br>0.3586 | 0.3800<br>0.3501 | 0.3745 | Ave | | 0.3700 | | | | 2.8 | | 15.0 | | | |
| Carbon disulfide | +++++<br>0.7777 | +++++<br>0.8035 | 0.6571<br>0.7823 | 0.6831<br>0.7752 | 0.7150 | Ave | | 0.7420 | | | 0.1000 | 7.6 | | 15.0 | | | |
| 3-Chloro-1-propene | 0.1133<br>0.1595 | 0.1186<br>0.1637 | 0.1545<br>0.1568 | 0.1443<br>0.1569 | 0.1528 | Ave | | 0.1467 | | | | 12.4 | | 15.0 | | | |
| Methyl acetate | 0.1968<br>0.1883 | 0.1858<br>0.1889 | 0.1900<br>0.1754 | 0.1789<br>0.1736 | 0.1960 | Ave | | 0.1860 | | | 0.1000 | 4.5 | | 15.0 | | | |
| Methylene Chloride | +++++<br>0.2820 | +++++<br>0.2745 | 0.4401<br>0.2582 | 0.3446<br>0.2521 | 0.3138 | Qua | 0.7353 | 0.2768 | -0.000111 | | 0.1000 | | | | 1.0000 | | 0.9900 |
| tert-Butyl alcohol (TBA) | 0.0268<br>0.0244 | 0.0261<br>0.0252 | 0.0256<br>0.0227 | 0.0232<br>0.0228 | 0.0259 | Ave | | 0.0247 | | | | 6.3 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.2875<br>0.3010 | 0.2780<br>0.3011 | 0.3055<br>0.2874 | 0.2940<br>0.2838 | 0.3088 | Ave | | 0.2941 | | | 0.1000 | 3.6 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.5907<br>0.6560 | 0.6124<br>0.6597 | 0.6455<br>0.6222 | 0.6055<br>0.6055 | 0.6553 | Ave | | 0.6281 | | | 0.1000 | 4.2 | | 15.0 | | | |
| Acrylonitrile | 0.1054<br>0.1049 | 0.1101<br>0.1046 | 0.1096<br>0.0948 | 0.1021<br>0.0930 | 0.1093 | Ave | | 0.1038 | | | | 6.0 | | 15.0 | | | |
| Hexane | 0.0913<br>0.0804 | 0.0651<br>0.0819 | 0.0896<br>0.0802 | 0.0799<br>0.0801 | 0.0814 | Ave | | 0.0811 | | | | 9.1 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.6065<br>0.6523 | 0.6496<br>0.6635 | 0.6625<br>0.6431 | 0.6244<br>0.6384 | 0.6558 | Ave | | 0.6440 | | | 0.1000 | 2.9 | | 15.0 | | | |
| Vinyl acetate | 0.0232<br>0.0341 | 0.0338<br>0.0370 | 0.0305<br>0.0354 | 0.0309<br>0.0364 | 0.0369 | Ave | | 0.0331 | | | | 13.4 | | 15.0 | | | |
| 2,2-Dichloropropane | 0.2668<br>0.3106 | 0.2490<br>0.3006 | 0.2757<br>0.3155 | 0.2904<br>0.2872 | 0.2925 | Ave | | 0.2876 | | | | 7.3 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.2639<br>0.3116 | 0.2890<br>0.3131 | 0.3133<br>0.3013 | 0.3126<br>0.2969 | 0.3112 | Ave | | 0.3014 | | | 0.1000 | 5.5 | | 15.0 | | | |
| 2-Butanone (MEK) | 0.1059<br>0.1103 | 0.1280<br>0.1145 | 0.1184<br>0.1087 | 0.1067<br>0.1077 | 0.1164 | Ave | | 0.1130 | | | 0.0500 | 6.4 | | 15.0 | | | |
| Bromochloromethane | 0.1211<br>0.1200 | 0.1279<br>0.1216 | 0.1177<br>0.1145 | 0.1173<br>0.1129 | 0.1240 | Ave | | 0.1197 | | | | 3.9 | | 15.0 | | | |
| Tetrahydrofuran | 0.0861<br>0.0732 | 0.0848<br>0.0753 | 0.0754<br>0.0695 | 0.0706<br>0.0688 | 0.0760 | Ave | | 0.0755 | | | | 8.2 | | 15.0 | | | |
| Chloroform | 0.4929<br>0.4944 | 0.5091<br>0.4997 | 0.4999<br>0.4837 | 0.4844<br>0.4787 | 0.5031 | Ave | | 0.4940 | | | | 2.0 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.3319<br>0.4423 | 0.3651<br>0.4534 | 0.4196<br>0.4375 | 0.4073<br>0.4350 | 0.4304 | Ave | | 0.4136 | | | 0.1000 | 9.7 | | 15.0 | | | |
| Cyclohexane | 0.6137<br>0.8524 | 0.6668<br>0.8401 | 0.8556<br>0.8002 | 0.8295<br>0.7872 | 0.8528 | Ave | | 0.7887 | | | 0.1000 | 11.2 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon tetrachloride | 0.2475<br>0.3648 | 0.2798<br>0.3753 | 0.3145<br>0.3681 | 0.3238<br>0.3653 | 0.3414 | Ave | | 0.3312 | | | 0.1000 | 13.4 | | 15.0 | | | |
| 1,1-Dichloropropene | 0.3223<br>0.3768 | 0.3284<br>0.3843 | 0.3741<br>0.3669 | 0.3744<br>0.3633 | 0.3884 | Ave | | 0.3643 | | | | 6.4 | | 15.0 | | | |
| Benzene | 1.0335<br>1.1238 | 1.0928<br>1.1312 | 1.1624<br>1.0841 | 1.1139<br>1.0648 | 1.1460 | Ave | | 1.1058 | | | 0.5000 | 3.7 | | 15.0 | | | |
| Isobutyl alcohol | 0.0084<br>0.0082 | 0.0085<br>0.0080 | 0.0094<br>0.0071 | 0.0077<br>0.0068 | 0.0087 | Ave | | 0.0081 | | | | 10.2 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.3941<br>0.4034 | 0.4149<br>0.4022 | 0.4116<br>0.3836 | 0.3912<br>0.3832 | 0.4179 | Ave | | 0.4002 | | | 0.1000 | 3.3 | | 15.0 | | | |
| n-Heptane | 0.0593<br>0.0818 | 0.0561<br>0.0840 | 0.0779<br>0.0812 | 0.0789<br>0.0793 | 0.0809 | Ave | | 0.0755 | | | | 13.6 | | 15.0 | | | |
| Trichloroethene | 0.2542<br>0.2825 | 0.2744<br>0.2855 | 0.2813<br>0.2763 | 0.2749<br>0.2710 | 0.2789 | Ave | | 0.2754 | | | 0.2000 | 3.3 | | 15.0 | | | |
| Methylcyclohexane | 0.3999<br>0.5630 | 0.4085<br>0.5667 | 0.5434<br>0.5379 | 0.5386<br>0.5259 | 0.5529 | Ave | | 0.5152 | | | 0.1000 | 12.5 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.2950<br>0.3233 | 0.2865<br>0.3257 | 0.3191<br>0.3207 | 0.3147<br>0.3177 | 0.3304 | Ave | | 0.3148 | | | 0.1000 | 4.6 | | 15.0 | | | |
| Dibromomethane | 0.1128<br>0.1268 | 0.1209<br>0.1299 | 0.1238<br>0.1240 | 0.1180<br>0.1234 | 0.1317 | Ave | | 0.1235 | | | | 4.7 | | 15.0 | | | |
| 1,4-Dioxane | 0.0019<br>0.0019 | 0.0020<br>0.0018 | 0.0019<br>0.0017 | 0.0018<br>0.0017 | 0.0020 | Ave | | 0.0019 | | | | 6.1 | | 15.0 | | | |
| Bromodichloromethane | 0.2163<br>0.3165 | 0.2426<br>0.3316 | 0.2765<br>0.3335 | 0.2644<br>0.3343 | 0.2928 | Ave | | 0.2898 | | | 0.1000 | 14.8 | | 15.0 | | | |
| 2-Chloroethyl vinyl ether | 0.0737<br>0.0999 | 0.0857<br>0.1065 | 0.0933<br>0.1041 | 0.0870<br>0.1073 | 0.0983 | Ave | | 0.0951 | | | | 11.8 | | 15.0 | | | |
| cis-1,3-Dichloropropene | +++++<br>0.3488 | +++++<br>0.3715 | 0.2701<br>0.3657 | 0.2919<br>0.3696 | 0.3296 | Ave | | 0.3353 | | | 0.1500 | 12.0 | | 15.0 | | | |
| 4-Methyl-2-pentanone (MIBK) | 0.3332<br>0.3715 | 0.3736<br>0.3858 | 0.3850<br>0.3499 | 0.3768<br>0.3448 | 0.3955 | Ave | | 0.3685 | | | 0.0500 | 5.7 | | 15.0 | | | |
| Toluene | 1.7560<br>1.8529 | 1.8581<br>1.9032 | 1.9298<br>1.8289 | 1.9095<br>1.7918 | 1.9412 | Ave | | 1.8635 | | | 0.4000 | 3.4 | | 15.0 | | | |
| trans-1,3-Dichloropropene | +++++<br>0.4199 | +++++<br>0.4568 | 0.3521<br>0.4527 | 0.3887<br>0.4582 | 0.4194 | Ave | | 0.4211 | | | 0.1000 | 9.4 | | 15.0 | | | |
| Ethyl methacrylate | 0.3108<br>0.3407 | 0.2940<br>0.3485 | 0.3045<br>0.3319 | 0.3067<br>0.3320 | 0.3538 | Ave | | 0.3247 | | | | 6.6 | | 15.0 | | | |
| 1,1,2-Trichloroethane | 0.3322<br>0.2812 | 0.2798<br>0.2877 | 0.3011<br>0.2722 | 0.2945<br>0.2702 | 0.2998 | Ave | | 0.2910 | | | 0.1000 | 6.6 | | 15.0 | | | |
| Tetrachloroethene | 0.3106<br>0.3371 | 0.3437<br>0.3518 | 0.3720<br>0.3355 | 0.3559<br>0.3312 | 0.3611 | Ave | | 0.3443 | | | 0.2000 | 5.3 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,3-Dichloropropane | 0.5096<br>0.4787 | 0.4843<br>0.4908 | 0.5023<br>0.4682 | 0.5029<br>0.4659 | 0.5075 | Ave | | 0.4900 | | | | 3.4 | | 15.0 | | | |
| 2-Hexanone | 0.2575<br>0.2403 | 0.2721<br>0.2417 | 0.2839<br>0.2214 | 0.2647<br>0.2183 | 0.2710 | Ave | | 0.2523 | | | 0.0500 | 9.2 | | 15.0 | | | |
| Dibromochloromethane | +++++<br>0.2828 | +++++<br>0.3185 | 0.2307<br>0.3152 | 0.2480<br>0.3131 | 0.2713 | Ave | | 0.2828 | | | | 12.3 | | 15.0 | | | |
| 1,2-Dibromoethane | 0.2578<br>0.2540 | 0.2468<br>0.2695 | 0.2545<br>0.2540 | 0.2476<br>0.2507 | 0.2699 | Ave | | 0.2561 | | | | 3.3 | | 15.0 | | | |
| Chlorobenzene | 1.1283<br>1.0422 | 1.1149<br>1.0849 | 1.1343<br>1.0176 | 1.1032<br>0.9905 | 1.0770 | Ave | | 1.0770 | | | 0.3000 | 4.7 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.2682<br>0.3744 | 0.3193<br>0.3887 | 0.3422<br>0.3659 | 0.3642<br>0.3571 | 0.3731 | Ave | | 0.3503 | | | | 10.5 | | 15.0 | | | |
| Ethylbenzene | 0.5951<br>0.6062 | 0.6488<br>0.6127 | 0.6416<br>0.5767 | 0.6241<br>0.5640 | 0.6134 | Ave | | 0.6092 | | | | 4.6 | | 15.0 | | | |
| m-Xylene & p-Xylene | 0.6293<br>0.7288 | 0.7435<br>0.7569 | 0.7919<br>0.7239 | 0.7625<br>0.7141 | 0.7585 | Ave | | 0.7344 | | | | 6.3 | | 15.0 | | | |
| o-Xylene | 0.6960<br>0.7398 | 0.6683<br>0.7498 | 0.7733<br>0.6927 | 0.7605<br>0.6753 | 0.7713 | Ave | | 0.7252 | | | | 5.8 | | 15.0 | | | |
| Styrene | 0.9671<br>1.1084 | 0.9701<br>1.1520 | 1.0851<br>1.0913 | 1.1030<br>1.0772 | 1.1652 | Ave | | 1.0799 | | | 0.3000 | 6.4 | | 15.0 | | | |
| Bromoform | 0.1197<br>0.1509 | 0.1386<br>0.1696 | 0.1262<br>0.1737 | 0.1259<br>0.1725 | 0.1435 | Ave | | 0.1467 | | | 0.1000 | 14.4 | | 15.0 | | | |
| Isopropylbenzene | 1.7485<br>2.0968 | 1.8947<br>2.1588 | 2.1170<br>2.0393 | 2.2006<br>1.9720 | 2.1982 | Ave | | 2.0473 | | | 0.1000 | 7.4 | | 15.0 | | | |
| Bromobenzene | 0.8099<br>0.8602 | 0.7750<br>0.8893 | 0.8982<br>0.8567 | 0.8290<br>0.8429 | 0.8798 | Ave | | 0.8490 | | | | 4.7 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 0.5506<br>0.6852 | 0.6400<br>0.6934 | 0.6674<br>0.6364 | 0.6224<br>0.6111 | 0.7209 | Ave | | 0.6475 | | | 0.3000 | 7.9 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.1211<br>0.1914 | 0.2024<br>0.2026 | 0.1913<br>0.1867 | 0.1926<br>0.1835 | 0.2057 | Ave | | 0.1864 | | | | 13.7 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | 0.2951<br>0.2807 | 0.2823<br>0.2935 | 0.2571<br>0.2837 | 0.2488<br>0.2858 | 0.2855 | Ave | | 0.2792 | | | | 5.6 | | 15.0 | | | |
| N-Propylbenzene | 0.8403<br>1.0899 | 0.8401<br>1.1354 | 1.0953<br>1.0951 | 1.0423<br>1.0667 | 1.1339 | Ave | | 1.0377 | | | | 11.1 | | 15.0 | | | |
| 2-Chlorotoluene | 0.7782<br>0.9360 | 0.7601<br>0.9575 | 0.8929<br>0.9090 | 0.9212<br>0.8801 | 0.9550 | Ave | | 0.8878 | | | | 8.1 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 2.6744<br>3.6914 | 2.7570<br>3.7764 | 3.3763<br>3.5538 | 3.3406<br>3.4306 | 3.6689 | Ave | | 3.3633 | | | | 11.8 | | 15.0 | | | |
| 4-Chlorotoluene | 0.8357<br>0.8953 | 0.8372<br>0.8987 | 0.9633<br>0.8755 | 0.8848<br>0.8685 | 0.9177 | Ave | | 0.8863 | | | | 4.5 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| tert-Butylbenzene | 2.2481<br>3.3321 | 2.5262<br>3.3872 | 3.0760<br>3.2252 | 3.0255<br>3.0788 | 3.3080 | Ave | | 3.0230 | | | | 12.8 | | 15.0 | | | |
| 1,2,4-Trimethylbenzene | 2.6774<br>3.6555 | 2.8117<br>3.7355 | 3.3662<br>3.5681 | 3.4053<br>3.4610 | 3.6683 | Ave | | 3.3721 | | | | 11.2 | | 15.0 | | | |
| sec-Butylbenzene | 3.2932<br>4.8364 | 3.6793<br>4.9728 | 4.4339<br>4.7070 | 4.5166<br>4.5183 | 4.7991 | Ave | | 4.4174 | | | | 12.8 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.6141<br>1.6288 | 1.6528<br>1.6255 | 1.7272<br>1.5617 | 1.6459<br>1.5144 | 1.6868 | Ave | | 1.6286 | | | 0.6000 | 3.9 | | 15.0 | | | |
| p-Isopropyltoluene | 2.8785<br>3.9630 | 3.0070<br>4.0561 | 3.8195<br>3.8320 | 3.7455<br>3.6576 | 3.9925 | Ave | | 3.6613 | | | | 11.7 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 1.7499<br>1.6007 | 1.6690<br>1.6056 | 1.7016<br>1.5732 | 1.6269<br>1.5403 | 1.6507 | Ave | | 1.6353 | | | 0.5000 | 4.0 | | 15.0 | | | |
| n-Butylbenzene | 2.8641<br>3.7743 | 2.9377<br>3.8919 | 3.6561<br>3.7690 | 3.5765<br>3.6431 | 3.8352 | Ave | | 3.5498 | | | | 10.7 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.7323<br>1.5627 | 1.6185<br>1.5512 | 1.7576<br>1.4688 | 1.5951<br>1.4159 | 1.6153 | Ave | | 1.5908 | | | 0.4000 | 6.9 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | 0.0965<br>0.1114 | 0.0893<br>0.1227 | 0.1176<br>0.1166 | 0.0839<br>0.1131 | 0.1080 | Ave | | 0.1066 | | | 0.0500 | 12.7 | | 15.0 | | | |
| 1,2,4-Trichlorobenzene | 1.2595<br>1.2579 | 1.2581<br>1.2747 | 1.3618<br>1.2022 | 1.2261<br>1.1559 | 1.3274 | Ave | | 1.2582 | | | 0.2000 | 4.9 | | 15.0 | | | |
| Hexachlorobutadiene | 0.6872<br>0.8507 | 0.8024<br>0.8827 | 0.8557<br>0.8148 | 0.8226<br>0.7684 | 0.8606 | Ave | | 0.8161 | | | | 7.3 | | 15.0 | | | |
| Naphthalene | 2.2789<br>2.5754 | 2.3038<br>2.5880 | 2.7275<br>2.4034 | 2.3883<br>2.2463 | 2.6721 | Ave | | 2.4649 | | | | 7.3 | | 15.0 | | | |
| 1,2,3-Trichlorobenzene | 1.2783<br>1.1998 | 1.2015<br>1.1854 | 1.3244<br>1.1044 | 1.1873<br>1.0128 | 1.2792 | Ave | | 1.1970 | | | | 7.9 | | 15.0 | | | |
| Dibromofluoromethane (Surr) | 0.2554<br>0.2127 | 0.2475<br>0.2245 | 0.2324<br>0.2169 | 0.2263<br>0.2198 | 0.2391 | Ave | | 0.2305 | | | | 6.3 | | 15.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | +++++<br>0.2812 | 0.3613<br>0.2947 | 0.3291<br>0.2912 | 0.3036<br>0.2990 | 0.3198 | Ave | | 0.3100 | | | | 8.3 | | 15.0 | | | |
| Toluene-d8 (Surr) | +++++<br>1.4561 | 1.7268<br>1.5560 | 1.6708<br>1.5304 | 1.6593<br>1.5592 | 1.7011 | Ave | | 1.6075 | | | | 5.9 | | 15.0 | | | |
| 4-Bromofluorobenzene (Surr) | +++++<br>1.0336 | +++++<br>1.1097 | 1.2422<br>1.1311 | 1.1090<br>1.1480 | 1.2017 | Ave | | 1.1393 | | | | 6.0 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-212606/11 | 147308.D |
| Level 2 | IC 240-212606/10 | 147307.D |
| Level 3 | IC 240-212606/9 | 147306.D |
| Level 4 | IC 240-212606/8 | 147305.D |
| Level 5 | IC 240-212606/7 | 147304.D |
| Level 6 | ICIS 240-212606/6 | 147303.D |
| Level 7 | IC 240-212606/5 | 147302.D |
| Level 8 | IC 240-212606/4 | 147301.D |
| Level 9 | IC 240-212606/3 | 147300.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Dichlorodifluoromethane | FB | Ave | 5871<br>419333 | 12661<br>826752 | 38481<br>1638337 | 79778<br>1967734 | 159816 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Chloromethane | FB | Ave | 11373<br>622955 | 20429<br>1252342 | 56747<br>2438111 | 117767<br>2802393 | 235327 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Vinyl chloride | FB | Ave | 7204<br>408074 | 13514<br>839111 | 39274<br>1617382 | 76856<br>1927305 | 155644 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Butadiene | FB | Ave | 6923<br>495753 | 14346<br>1003660 | 44709<br>1963149 | 97227<br>2345207 | 188492 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Bromomethane | FB | Ave | 3716<br>156173 | 6947<br>+++++ | 16327<br>+++++ | 34300<br>+++++ | 65725 | 1.00<br>50.0 | 2.00<br>+++++ | 5.00<br>+++++ | 10.0<br>+++++ | 20.0 |
| Chloroethane | FB | Ave | 4856<br>209466 | 6965<br>394781 | 21010<br>+++++ | 42227<br>+++++ | 81510 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>+++++ | 10.0<br>+++++ | 20.0 |
| Dichlorofluoromethane | FB | Ave | 8681<br>444994 | 15322<br>854115 | 39828<br>1737417 | 89907<br>2044076 | 161421 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Trichlorofluoromethane | FB | Ave | 7820<br>469944 | 13626<br>937500 | 40109<br>1823298 | 88470<br>2211446 | 177978 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl ether | FB | Ave | +++++<br>283284 | +++++<br>561509 | 29032<br>1064646 | 54103<br>1278520 | 113841 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Acrolein | FB | Lin1 | 5499<br>218567 | 9009<br>452990 | 34458<br>810395 | 44584<br>1004148 | 91896 | 5.00<br>250 | 10.0<br>500 | 25.0<br>1000 | 50.0<br>1250 | 100 |
| 1,1-Dichloroethene | FB | Ave | 4586<br>323142 | 11115<br>684544 | 34169<br>1275244 | 64535<br>1490519 | 132072 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroetha ne | FB | Ave | 3913<br>276872 | 9251<br>568070 | 28361<br>1074115 | 51871<br>1281941 | 112563 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Acetone | FB | Qua | +++++<br>221530 | +++++<br>446091 | 34991<br>813218 | 53362<br>941400 | 102774 | +++++<br>100 | +++++<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Iodomethane | FB | Ave | 8786<br>488418 | 17673<br>959263 | 46310<br>1829838 | 94594<br>2186981 | 184371 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Carbon disulfide | FB | Ave | +++++<br>996181 | +++++<br>2076690 | 81978<br>3992337 | 170054<br>4842941 | 352045 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 3-Chloro-1-propene | FB | Ave | 2638<br>204278 | 5728<br>423070 | 19271<br>800139 | 35913<br>979966 | 75255 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methyl acetate | FB | Ave | 22906<br>1206297 | 44853<br>2441019 | 118518<br>4476273 | 222650<br>5422320 | 482624 | 5.00<br>250 | 10.0<br>500 | 25.0<br>1000 | 50.0<br>1250 | 100 |
| Methylene Chloride | FB | Qua | +++++<br>361235 | +++++<br>709387 | 54903<br>1317748 | 85786<br>1574955 | 154528 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| tert-Butyl alcohol (TBA) | FB | Ave | 6246<br>312962 | 12610<br>652025 | 31980<br>1156473 | 57745<br>1423418 | 127348 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| trans-1,2-Dichloroethene | FB | Ave | 6694<br>385522 | 13425<br>778312 | 38117<br>1466571 | 73178<br>1772943 | 152040 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methyl tert-butyl ether | FB | Ave | 13753<br>840271 | 29573<br>1705065 | 80533<br>3175146 | 150726<br>3782570 | 322668 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Acrylonitrile | FB | Ave | 24531<br>1343873 | 53178<br>2704516 | 136763<br>4840214 | 254223<br>5807759 | 537940 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| Hexane | FB | Ave | 2125<br>102982 | 3142<br>211731 | 11185<br>409313 | 19878<br>500287 | 40067 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1-Dichloroethane | FB | Ave | 14120<br>835497 | 31368<br>1714925 | 82654<br>3281739 | 155431<br>3988543 | 322893 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Vinyl acetate | FB | Ave | 540<br>43742 | 1633<br>95718 | 3804<br>180417 | 7684<br>227375 | 18152 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2,2-Dichloropropane | FB | Ave | 6212<br>397873 | 12025<br>777050 | 34402<br>1609897 | 72287<br>1794262 | 144041 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| cis-1,2-Dichloroethene | FB | Ave | 6143<br>399172 | 13958<br>809127 | 39088<br>1537416 | 77811<br>1854690 | 153236 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Butanone (MEK) | FB | Ave | 4932<br>282507 | 12367<br>592067 | 29540<br>1109488 | 53122<br>1345636 | 114610 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Bromochloromethane | FB | Ave | 2819<br>153736 | 6177<br>314240 | 14681<br>584364 | 29199<br>705188 | 61039 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Tetrahydrofuran | FB | Ave | 4010<br>187593 | 8188<br>389332 | 18808<br>709841 | 35143<br>859859 | 74810 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Chloroform | FB | Ave | 11476<br>633312 | 24584<br>1291586 | 62366<br>2468431 | 120584<br>2990336 | 247732 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,1-Trichloroethane | FB | Ave | 7728<br>566525 | 17632<br>1171935 | 52352<br>2232718 | 101392<br>2717739 | 211942 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Cyclohexane | FB | Ave | 14287<br>1091809 | 32199<br>2171202 | 106745<br>4083646 | 206503<br>4918071 | 419919 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Carbon tetrachloride | FB | Ave | 5762<br>467218 | 13511<br>969904 | 39240<br>1878751 | 80601<br>2282250 | 168076 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1-Dichloropropene | FB | Ave | 7503<br>482629 | 15860<br>993188 | 46673<br>1872225 | 93194<br>2269592 | 191238 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Benzene | FB | Ave | 24060<br>1439453 | 52774<br>2923674 | 145026<br>5532307 | 277304<br>6652146 | 564247 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Isobutyl alcohol | FB | Ave | 4904<br>263907 | 10301<br>518655 | 29415<br>906552 | 47705<br>1059248 | 107668 | 25.0<br>1250 | 50.0<br>2500 | 125<br>5000 | 250<br>6250 | 500 |
| 1,2-Dichloroethane | FB | Ave | 9175<br>516699 | 20035<br>1039532 | 51349<br>1957435 | 97389<br>2394295 | 205784 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| n-Heptane | FB | Ave | 1380<br>104720 | 2709<br>217129 | 9718<br>414474 | 19639<br>495570 | 39816 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Trichloroethene | FB | Ave | 5917<br>361820 | 13249<br>737799 | 35092<br>1409933 | 68425<br>1692775 | 137337 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methylcyclohexane | FB | Ave | 9310<br>721138 | 19728<br>1464792 | 67792<br>2744795 | 134079<br>3285514 | 272228 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dichloropropane | FB | Ave | 6867<br>414151 | 13834<br>841807 | 39813<br>1636687 | 78347<br>1984906 | 162657 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Dibromomethane | FB | Ave | 2626<br>162359 | 5840<br>335689 | 15440<br>632675 | 29368<br>771219 | 64857 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,4-Dioxane | FB | Ave | 887<br>48044 | 1957<br>94875 | 4664<br>176535 | 8933<br>215441 | 20175 | 20.0<br>1000 | 40.0<br>2000 | 100<br>4000 | 200<br>5000 | 400 |
| Bromodichloromethane | FB | Ave | 5035<br>405424 | 11715<br>857006 | 34493<br>1701789 | 65817<br>2088773 | 144174 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Chloroethyl vinyl ether | FB | Ave | 3433<br>256022 | 8273<br>550543 | 23278<br>1062923 | 43328<br>1340964 | 96811 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| cis-1,3-Dichloropropene | FB | Ave | +++++<br>446732 | +++++<br>960279 | 33701<br>1866444 | 72658<br>2309301 | 162293 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | CBZ | Ave | 8896<br>584970 | 20976<br>1198696 | 55579<br>2174185 | 107347<br>2647060 | 229049 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Toluene | CBZ | Ave | 23439<br>1458768 | 52162<br>2956356 | 139303<br>5682435 | 272022<br>6877637 | 562045 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| trans-1,3-Dichloropropene | CBZ | Ave | +++++<br>330594 | +++++<br>709556 | 25419<br>1406473 | 55379<br>1758690 | 121435 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl methacrylate | CBZ | Ave | 4149<br>268222 | 8252<br>541319 | 21977<br>1031104 | 43685<br>1274320 | 102425 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,2-Trichloroethane | CBZ | Ave | 4434<br>221353 | 7855<br>446892 | 21734<br>845653 | 41957<br>1037101 | 86815 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Tetrachloroethene | CBZ | Ave | 4146<br>265409 | 9650<br>546491 | 26856<br>1042254 | 50698<br>1271281 | 104563 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3-Dichloropropane | CBZ | Ave | 6802<br>376892 | 13597<br>762406 | 36259<br>1454663 | 71647<br>1788276 | 146941 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Hexanone | CBZ | Ave | 6873<br>378399 | 15278<br>750989 | 40986<br>1375916 | 75413<br>1675849 | 156916 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Dibromochloromethane | CBZ | Ave | +++++<br>222618 | +++++<br>494736 | 16654<br>979386 | 35323<br>1201593 | 78552 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1,2-Dibromoethane | CBZ | Ave | 3441<br>199936 | 6929<br>418627 | 18368<br>789296 | 35270<br>962215 | 78139 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Chlorobenzene | CBZ | Ave | 15061<br>820507 | 31298<br>1685148 | 81879<br>3161618 | 157153<br>3801868 | 311842 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,1,2-Tetrachloroethane | CBZ | Ave | 3580<br>294759 | 8963<br>603739 | 24702<br>1136904 | 51880<br>1370710 | 108027 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethylbenzene | CBZ | Ave | 7943<br>477219 | 18214<br>951746 | 46312<br>1791696 | 88903<br>2164935 | 177597 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| m-Xylene & p-Xylene | CBZ | Ave | 8400<br>573737 | 20873<br>1175688 | 57167<br>2249090 | 108629<br>2741018 | 219601 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| o-Xylene | CBZ | Ave | 9290<br>582431 | 18762<br>1164691 | 55819<br>2152239 | 108337<br>2591902 | 223316 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Styrene | CBZ | Ave | 12909<br>872604 | 27233<br>1789449 | 78330<br>3390778 | 157132<br>4134767 | 337371 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Bromoform | CBZ | Ave | 1598<br>118785 | 3892<br>263387 | 9107<br>539628 | 17930<br>662011 | 41561 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Isopropylbenzene | CBZ | Ave | 23339<br>1650732 | 53189<br>3353369 | 152820<br>6336102 | 313488<br>7569173 | 636450 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Bromobenzene | DCB | Ave | 5777<br>322109 | 12037<br>654908 | 33615<br>1238302 | 63294<br>1519489 | 128751 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,1,2,2-Tetrachloroethane | DCB | Ave | 3927<br>256576 | 9940<br>510672 | 24976<br>919809 | 47520<br>1101627 | 105495 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,3-Trichloropropane | DCB | Ave | 864<br>71686 | 3144<br>149211 | 7158<br>269890 | 14703<br>330871 | 30098 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| trans-1,4-Dichloro-2-butene | DCB | Ave | 2105<br>105103 | 4384<br>216157 | 9621<br>410063 | 18997<br>515186 | 41779 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| N-Propylbenzene | DCB | Ave | 5994<br>408139 | 13049<br>836161 | 40993<br>1582897 | 79579<br>1923014 | 165943 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Chlorotoluene | DCB | Ave | 5551<br>350484 | 11806<br>705190 | 33416<br>1313804 | 70336<br>1586630 | 139760 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3,5-Trimethylbenzene | DCB | Ave | 19076<br>1382278 | 42822<br>2781225 | 126357<br>5136692 | 255059<br>6184624 | 536911 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 4-Chlorotoluene | DCB | Ave | 5961<br>335245 | 13003<br>661900 | 36050<br>1265427 | 67552<br>1565650 | 134298 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| tert-Butylbenzene | DCB | Ave | 16035<br>1247755 | 39237<br>2494554 | 115120<br>4661764 | 231001<br>5550377 | 484092 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,4-Trimethylbenzene | DCB | Ave | 19097<br>1368854 | 43671<br>2751107 | 125980<br>5157358 | 259998<br>6239391 | 536824 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| sec-Butylbenzene | DCB | Ave | 23490<br>1811021 | 57147<br>3662298 | 165940<br>6803540 | 344844<br>8145446 | 702303 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3-Dichlorobenzene | DCB | Ave | 11513<br>609919 | 25671<br>1197171 | 64639<br>2257263 | 125668<br>2730042 | 246844 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 12:29 Calibration End Date: 12/30/2015 15:21 Calibration ID: 32237

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| p-Isopropyltoluene | DCB | Ave | 20532<br>1483977 | 46705<br>2987227 | 142946<br>5538742 | 285971<br>6593837 | 584270 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,4-Dichlorobenzene | DCB | Ave | 12482<br>599393 | 25922<br>1182499 | 63682<br>2273928 | 124212<br>2776758 | 241560 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| n-Butylbenzene | DCB | Ave | 20429<br>1413342 | 45628<br>2866292 | 136830<br>5447690 | 273066<br>6567734 | 561253 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dichlorobenzene | DCB | Ave | 12356<br>585151 | 25139<br>1142395 | 65779<br>2122980 | 121785<br>2552468 | 236378 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | DCB | Ave | 688<br>41703 | 1387<br>90331 | 4403<br>168526 | 6403<br>203958 | 15806 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,4-Trichlorobenzene | DCB | Ave | 8984<br>471040 | 19541<br>938771 | 50966<br>1737656 | 93617<br>2083801 | 194259 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Hexachlorobutadiene | DCB | Ave | 4902<br>318558 | 12463<br>650106 | 32023<br>1177742 | 62807<br>1385206 | 125945 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Naphthalene | DCB | Ave | 16255<br>964386 | 35782<br>1905999 | 102076<br>3473895 | 182350<br>4049536 | 391041 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2,3-Trichlorobenzene | DCB | Ave | 9118<br>449280 | 18661<br>873029 | 49567<br>1596336 | 90651<br>1825837 | 187196 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Dibromofluoromethane (Surr) | FB | Ave | 5947<br>272413 | 11951<br>580216 | 28996<br>1106784 | 56336<br>1373048 | 117734 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Ave | +++++<br>360241 | 17448<br>761644 | 41065<br>1486058 | 75587<br>1867667 | 157455 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Toluene-d8 (Surr) | CBZ | Ave | +++++<br>1146345 | 48475<br>2416949 | 120605<br>4755110 | 236385<br>5984676 | 492531 | +++++<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 4-Bromofluorobenzene (Surr) | DCB | Ave | +++++<br>387052 | +++++<br>817275 | 46489<br>1634906 | 84672<br>2069550 | 175852 | +++++<br>50.0 | +++++<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

Curve Type Legend:

Ave = Average ISTD
Lin1 = Linear 1/conc ISTD
Qua = Quadratic ISTD

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-59482-1  Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 15:42  Calibration End Date: 12/30/2015 18:34  Calibration ID: 32239

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-212606/20 | 147317.D |
| Level 2 | IC 240-212606/19 | 147316.D |
| Level 3 | IC 240-212606/18 | 147315.D |
| Level 4 | IC 240-212606/17 | 147314.D |
| Level 5 | IC 240-212606/16 | 147313.D |
| Level 6 | IC 240-212606/15 | 147312.D |
| Level 7 | IC 240-212606/14 | 147311.D |
| Level 8 | IC 240-212606/13 | 147310.D |
| Level 9 | IC 240-212606/12 | 147309.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 LVL 7 | LVL 3 LVL 8 | LVL 4 LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Acetonitrile | 0.0354 0.0296 | 0.0354 0.0329 | 0.0304 0.0298 | 0.0312 0.0320 | 0.0308 | Ave | | 0.0319 | | | | 6.9 | | 15.0 | | | |
| 2-Chloro-1,3-butadiene | 0.5267 0.7036 | 0.5241 0.7474 | 0.6123 0.7356 | 0.6721 0.7515 | 0.6963 | Ave | | 0.6633 | | | | 13.4 | | 15.0 | | | |
| Diisopropyl ether | 0.1886 0.2334 | 0.2163 0.2442 | 0.2171 0.2316 | 0.2141 0.2371 | 0.2274 | Ave | | 0.2233 | | | | 7.4 | | 15.0 | | | |
| Ethyl t-butyl ether | 0.8514 1.0598 | 0.9172 1.1400 | 0.9049 1.0975 | 0.9690 1.1357 | 1.0429 | Ave | | 1.0131 | | | | 10.5 | | 15.0 | | | |
| Ethyl acetate | 0.2100 0.2184 | 0.1888 0.2218 | 0.1959 0.2021 | 0.2065 0.2188 | 0.2169 | Ave | | 0.2088 | | | | 5.5 | | 15.0 | | | |
| Propionitrile | 0.0383 0.0379 | 0.0401 0.0401 | 0.0366 0.0363 | 0.0383 0.0388 | 0.0396 | Ave | | 0.0384 | | | | 3.6 | | 15.0 | | | |
| Methacrylonitrile | 0.1296 0.1478 | 0.1453 0.1493 | 0.1390 0.1328 | 0.1503 0.1384 | 0.1516 | Ave | | 0.1427 | | | | 5.6 | | 15.0 | | | |
| Tert-amyl methyl ether | 0.5156 0.6672 | 0.5498 0.7269 | 0.5868 0.6865 | 0.6257 0.7186 | 0.6723 | Ave | | 0.6388 | | | | 11.7 | | 15.0 | | | |
| n-Butanol | 0.0103 0.0119 | 0.0115 0.0140 | 0.0091 0.0126 | 0.0111 0.0142 | 0.0114 | Ave | | 0.0118 | | | | 13.9 | | 15.0 | | | |
| Ethyl acrylate | 0.2546 0.3181 | 0.2502 0.3110 | 0.2735 0.2870 | 0.2734 0.3103 | 0.2902 | Ave | | 0.2854 | | | | 8.6 | | 15.0 | | | |
| Methyl methacrylate | 0.1662 0.2064 | 0.1707 0.2026 | 0.1698 0.1828 | 0.1816 0.1978 | 0.1961 | Ave | | 0.1860 | | | | 8.2 | | 15.0 | | | |
| 2-Nitropropane | 0.0416 0.0370 | 0.0343 0.0425 | 0.0325 0.0415 | 0.0337 0.0469 | 0.0317 | Ave | | 0.0380 | | | | 14.1 | | 15.0 | | | |
| n-Butyl acetate | 0.7901 0.6095 | 0.7350 0.6036 | 0.5430 0.5286 | 0.6001 0.6008 | 0.5924 | Ave | | 0.6226 | | | | 13.7 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 15:42 Calibration End Date: 12/30/2015 18:34 Calibration ID: 32239

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1-Chlorohexane | 0.4939<br>0.5358 | 0.4219<br>0.5405 | 0.4649<br>0.5445 | 0.4910<br>0.5595 | 0.5060 | Ave | | 0.5064 | | | | 8.7 | | 15.0 | | | |
| Cyclohexanone | 0.0108<br>0.0147 | 0.0127<br>0.0163 | 0.0140<br>0.0149 | 0.0139<br>0.0163 | 0.0140 | Ave | | 0.0142 | | | | 12.1 | | 15.0 | | | |
| 1,2,3-Trimethylbenzene | 2.6245<br>3.4683 | 2.8454<br>3.6539 | 2.9931<br>3.6054 | 3.2488<br>3.6720 | 3.3975 | Ave | | 3.2788 | | | | 11.6 | | 15.0 | | | |
| 1,3,5-Trichlorobenzene | 1.2285<br>1.2807 | 1.2778<br>1.3331 | 1.2687<br>1.3033 | 1.2897<br>1.3303 | 1.2961 | Ave | | 1.2898 | | | | 2.5 | | 15.0 | | | |
| 2-Methylnaphthalene | 1.3391<br>1.5219 | 1.2902<br>1.6776 | 1.2653<br>1.5531 | 1.3414<br>1.5514 | 1.6128 | Ave | | 1.4614 | | | | 10.5 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 15:42 Calibration End Date: 12/30/2015 18:34 Calibration ID: 32239

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-212606/20 | 147317.D |
| Level 2 | IC 240-212606/19 | 147316.D |
| Level 3 | IC 240-212606/18 | 147315.D |
| Level 4 | IC 240-212606/17 | 147314.D |
| Level 5 | IC 240-212606/16 | 147313.D |
| Level 6 | IC 240-212606/15 | 147312.D |
| Level 7 | IC 240-212606/14 | 147311.D |
| Level 8 | IC 240-212606/13 | 147310.D |
| Level 9 | IC 240-212606/12 | 147309.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Acetonitrile | FB | Ave | 8050<br>381248 | 16691<br>836473 | 35915<br>1479810 | 75285<br>1973898 | 148856 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| 2-Chloro-1,3-butadiene | FB | Ave | 11982<br>906998 | 24713<br>1898639 | 72396<br>3648912 | 162396<br>4629161 | 336600 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Diisopropyl ether | FB | Ave | 4290<br>300842 | 10201<br>620244 | 25668<br>1148932 | 51735<br>1460404 | 109917 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl t-butyl ether | FB | Ave | 19370<br>1366076 | 43245<br>2895796 | 106999<br>5444211 | 234143<br>6995373 | 504138 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Ethyl acetate | FB | Ave | 9555<br>562963 | 17808<br>1126679 | 46337<br>2004852 | 99804<br>2695800 | 209664 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| Propionitrile | FB | Ave | 8705<br>488992 | 18897<br>1017464 | 43264<br>1799603 | 92502<br>2392179 | 191650 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| Methacrylonitrile | FB | Ave | 29482<br>1904850 | 68497<br>3791818 | 164358<br>6587252 | 363218<br>8522886 | 732861 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| Tert-amyl methyl ether | FB | Ave | 11729<br>860017 | 25922<br>1846587 | 69383<br>3405251 | 151185<br>4426300 | 325028 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| n-Butanol | CBZ | Ave | 3470<br>227089 | 7855<br>507914 | 16280<br>901699 | 40119<br>1215606 | 83382 | 25.0<br>1250 | 50.0<br>2500 | 125<br>5000 | 250<br>6250 | 500 |
| Ethyl acrylate | FB | Ave | 5793<br>410014 | 11799<br>789971 | 32341<br>1423710 | 66058<br>1911099 | 140269 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| Methyl methacrylate | FB | Ave | 7562<br>532024 | 16100<br>1029437 | 40143<br>1813161 | 87763<br>2437020 | 189638 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| 2-Nitropropane | FB | Ave | 1895<br>95300 | 3236<br>215937 | 7678<br>411528 | 16302<br>577353 | 30610 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |
| n-Butyl acetate | CBZ | Ave | 10622<br>466530 | 20065<br>875681 | 38997<br>1509978 | 86776<br>2063646 | 172665 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1-Chlorohexane | CBZ | Ave | 6640<br>410111 | 11518<br>784090 | 33390<br>1555327 | 71000<br>1921563 | 147466 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 212606

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 12/30/2015 15:42 Calibration End Date: 12/30/2015 18:34 Calibration ID: 32239

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| Cyclohexanone | CBZ | Ave | 1452<br>112471 | 3470<br>236900 | 10056<br>425608 | 20164<br>560080 | 40836 | 10.0<br>500 | 20.0<br>1000 | 50.0<br>2000 | 100<br>2500 | 200 |
| 1,2,3-Trimethylbenzene | DCB | Ave | 18478<br>1388605 | 42188<br>2771667 | 109005<br>5188835 | 243094<br>6572001 | 504802 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 1,3,5-Trichlorobenzene | DCB | Ave | 8649<br>512752 | 18946<br>1011248 | 46203<br>1875695 | 96502<br>2380864 | 192581 | 1.00<br>50.0 | 2.00<br>100 | 5.00<br>200 | 10.0<br>250 | 20.0 |
| 2-Methylnaphthalene | DCB | Ave | 18856<br>1218632 | 38259<br>2545141 | 92161<br>4470426 | 200735<br>5553255 | 479256 | 2.00<br>100 | 4.00<br>200 | 10.0<br>400 | 20.0<br>500 | 40.0 |

Curve Type Legend:
Ave = Average ISTD

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-210738/7 | UXC5227.D |
| Level 2 | STD8260 240-210738/6 | UXC5226.D |
| Level 3 | STD8260 240-210738/5 | UXC5225.D |
| Level 4 | STD8260 240-210738/4 | UXC5224.D |
| Level 5 | STD8260 240-210738/3 | UXC5223.D |
| Level 6 | STD8260 240-210738/2 | UXC5222A.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | 0.4163<br>0.4078 | 0.4068 | 0.4254 | 0.3707 | 0.4170 | Ave | | 0.4073 | | | 0.1000 | 4.7 | | 15.0 | | | |
| Chloromethane | 0.5219<br>0.4236 | 0.4968 | 0.4810 | 0.4520 | 0.4482 | Ave | | 0.4706 | | | 0.1000 | 7.7 | | 15.0 | | | |
| Vinyl chloride | 0.4238<br>0.3959 | 0.4254 | 0.4328 | 0.3816 | 0.4118 | Ave | | 0.4119 | | | 0.1000 | 4.8 | | 15.0 | | | |
| Butadiene | 0.4000<br>0.3729 | 0.4010 | 0.4057 | 0.3545 | 0.3886 | Ave | | 0.3871 | | | | 5.1 | | 15.0 | | | |
| Bromomethane | 0.1520<br>0.1276 | 0.1359 | 0.1632 | 0.1052 | 0.1508 | Ave | | 0.1391 | | | 0.0500 | 15.0 | | 15.0 | | | |
| Chloroethane | 0.2483<br>0.1975 | 0.2298 | 0.2410 | 0.2016 | 0.2161 | Ave | | 0.2224 | | | | 9.4 | | 15.0 | | | |
| Dichlorofluoromethane | 0.5356<br>0.4476 | 0.4781 | 0.5225 | 0.4286 | 0.4896 | Ave | | 0.4837 | | | | 8.6 | | 15.0 | | | |
| Trichlorofluoromethane | 0.3464<br>0.3452 | 0.3375 | 0.3656 | 0.3031 | 0.3607 | Ave | | 0.3431 | | | 0.1000 | 6.5 | | 15.0 | | | |
| Ethyl ether | 0.3294<br>0.2319 | 0.2659 | 0.2650 | 0.2616 | 0.2453 | Ave | | 0.2665 | | | | 12.6 | | 15.0 | | | |
| Acrolein | 0.0386<br>0.0382 | 0.0359 | 0.0401 | 0.0381 | 0.0375 | Ave | | 0.0381 | | | | 3.6 | | 15.0 | | | |
| 1,1-Dichloroethene | 0.3301<br>0.3233 | 0.3228 | 0.3278 | 0.3221 | 0.3357 | Ave | | 0.3270 | | | 0.1000 | 1.6 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.2278<br>0.2524 | 0.2308 | 0.2386 | 0.2457 | 0.2660 | Ave | | 0.2436 | | | 0.0500 | 5.9 | | 15.0 | | | |
| Acetone | 0.1398<br>0.0712 | 0.1008 | 0.0883 | 0.0756 | 0.0699 | Lin1 | 0.1424 | 0.0687 | | | 0.0500 | | | | 0.9990 | | 0.9900 |
| Iodomethane | 0.5722<br>0.5292 | 0.5382 | 0.5474 | 0.5241 | 0.5408 | Ave | | 0.5420 | | | | 3.1 | | 15.0 | | | |
| Carbon disulfide | 0.9853<br>1.1135 | 0.9961 | 1.0372 | 1.0389 | 1.1172 | Ave | | 1.0481 | | | 0.1000 | 5.4 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 3-Chloro-1-propene | 0.2048<br>0.2099 | 0.1925 | 0.2008 | 0.2062 | 0.2137 | Ave | | 0.2047 | | | | 3.6 | | 15.0 | | | |
| Methyl acetate | 0.1922<br>0.1745 | 0.1844 | 0.1804 | 0.1688 | 0.1756 | Ave | | 0.1793 | | | 0.1000 | 4.6 | | 15.0 | | | |
| Methylene Chloride | 1.0108<br>0.3579 | 0.6706 | 0.5050 | 0.4172 | 0.3858 | Lin1 | 0.6767 | 0.3464 | | | 0.1000 | | | | 0.9990 | | 0.9900 |
| tert-Butyl alcohol (TBA) | 0.0168<br>0.0146 | 0.0161 | 0.0157 | 0.0153 | 0.0160 | Ave | | 0.0157 | | | | 4.8 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.8328<br>0.7862 | 0.7722 | 0.7967 | 0.7656 | 0.7965 | Ave | | 0.7917 | | | 0.1000 | 3.0 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.3711<br>0.3427 | 0.3572 | 0.3596 | 0.3445 | 0.3576 | Ave | | 0.3554 | | | 0.1000 | 3.0 | | 15.0 | | | |
| Acrylonitrile | 0.1004<br>0.0960 | 0.0954 | 0.0969 | 0.0926 | 0.0965 | Ave | | 0.0963 | | | | 2.6 | | 15.0 | | | |
| Hexane | 0.0824<br>0.0848 | 0.0812 | 0.0855 | 0.0874 | 0.0912 | Ave | | 0.0854 | | | | 4.2 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.6827<br>0.6354 | 0.6415 | 0.6713 | 0.6324 | 0.6534 | Ave | | 0.6528 | | | 0.1000 | 3.1 | | 15.0 | | | |
| Vinyl acetate | 0.4795<br>0.5103 | 0.4527 | 0.4862 | 0.4761 | 0.5072 | Ave | | 0.4853 | | | | 4.4 | | 15.0 | | | |
| 2,2-Dichloropropane | 0.3501<br>0.3522 | 0.3590 | 0.3702 | 0.3635 | 0.3799 | Ave | | 0.3625 | | | | 3.1 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.4002<br>0.3630 | 0.3771 | 0.3891 | 0.3682 | 0.3757 | Ave | | 0.3789 | | | 0.1000 | 3.6 | | 15.0 | | | |
| 2-Butanone (MEK) | 0.1134<br>0.0994 | 0.0969 | 0.1037 | 0.0961 | 0.0986 | Ave | | 0.1014 | | | 0.0500 | 6.4 | | 15.0 | | | |
| Bromochloromethane | 0.1664<br>0.1630 | 0.1684 | 0.1677 | 0.1636 | 0.1661 | Ave | | 0.1659 | | | | 1.3 | | 15.0 | | | |
| Tetrahydrofuran | 0.0948<br>0.0647 | 0.0813 | 0.0718 | 0.0642 | 0.0659 | Lin1 | 0.0642 | 0.0638 | | | | | | | 1.0000 | | 0.9900 |
| Chloroform | 0.6293<br>0.5814 | 0.5856 | 0.6053 | 0.5774 | 0.5921 | Ave | | 0.5952 | | | | 3.3 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.3918<br>0.4355 | 0.4209 | 0.4396 | 0.4299 | 0.4627 | Ave | | 0.4301 | | | 0.1000 | 5.4 | | 15.0 | | | |
| Cyclohexane | 0.6081<br>0.6300 | 0.5759 | 0.6142 | 0.6274 | 0.6608 | Ave | | 0.6194 | | | 0.1000 | 4.5 | | 15.0 | | | |
| 1,1-Dichloropropene | 0.4685<br>0.4667 | 0.4621 | 0.4795 | 0.4631 | 0.4854 | Ave | | 0.4709 | | | | 2.0 | | 15.0 | | | |
| Carbon tetrachloride | 0.3249<br>0.3889 | 0.3230 | 0.3466 | 0.3580 | 0.3938 | Ave | | 0.3559 | | | 0.1000 | 8.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Isobutyl alcohol | 0.0067<br>0.0059 | 0.0066 | 0.0062 | 0.0058 | 0.0062 | Ave | | 0.0062 | | | | 6.1 | | 15.0 | | | |
| Benzene | 1.6483<br>1.5004 | 1.5097 | 1.5340 | 1.4513 | 1.5265 | Ave | | 1.5284 | | | 0.5000 | 4.3 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.4708<br>0.4260 | 0.4438 | 0.4544 | 0.4273 | 0.4396 | Ave | | 0.4436 | | | 0.1000 | 3.8 | | 15.0 | | | |
| n-Heptane | 0.0878<br>0.0787 | 0.0700 | 0.0747 | 0.0799 | 0.0852 | Ave | | 0.0794 | | | | 8.3 | | 15.0 | | | |
| Trichloroethene | 0.3714<br>0.3510 | 0.3554 | 0.3681 | 0.3500 | 0.3641 | Ave | | 0.3600 | | | 0.2000 | 2.5 | | 15.0 | | | |
| Methylcyclohexane | 0.4769<br>0.5320 | 0.4925 | 0.5140 | 0.5315 | 0.5690 | Ave | | 0.5193 | | | 0.1000 | 6.3 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.3557<br>0.3473 | 0.3538 | 0.3550 | 0.3439 | 0.3510 | Ave | | 0.3511 | | | 0.1000 | 1.3 | | 15.0 | | | |
| Dibromomethane | 0.1801<br>0.1690 | 0.1649 | 0.1700 | 0.1625 | 0.1737 | Ave | | 0.1700 | | | | 3.7 | | 15.0 | | | |
| 1,4-Dioxane | 0.0027<br>0.0024 | 0.0025 | 0.0025 | 0.0025 | 0.0027 | Ave | | 0.0025 | | | | 5.1 | | 15.0 | | | |
| Bromodichloromethane | 0.3349<br>0.3950 | 0.3187 | 0.3573 | 0.3614 | 0.3916 | Ave | | 0.3598 | | | 0.1000 | 8.4 | | 15.0 | | | |
| 2-Chloroethyl vinyl ether | 0.1671<br>0.1770 | 0.1626 | 0.1732 | 0.1705 | 0.1793 | Ave | | 0.1716 | | | | 3.6 | | 15.0 | | | |
| cis-1,3-Dichloropropene | 0.4034<br>0.5079 | 0.4174 | 0.4549 | 0.4596 | 0.5021 | Ave | | 0.4575 | | | 0.1500 | 9.3 | | 15.0 | | | |
| 4-Methyl-2-pentanone (MIBK) | 0.2102<br>0.2060 | 0.1884 | 0.1953 | 0.1939 | 0.2029 | Ave | | 0.1994 | | | 0.0500 | 4.1 | | 15.0 | | | |
| Toluene | 2.0926<br>2.0094 | 2.0297 | 2.0385 | 1.9542 | 2.0306 | Ave | | 2.0258 | | | 0.4000 | 2.2 | | 15.0 | | | |
| trans-1,3-Dichloropropene | 0.3915<br>0.5343 | 0.4156 | 0.4712 | 0.4759 | 0.5219 | Ave | | 0.4684 | | | 0.1000 | 12.1 | | 15.0 | | | |
| Ethyl methacrylate | 0.3753<br>0.4447 | 0.3829 | 0.4186 | 0.4239 | 0.4409 | Ave | | 0.4144 | | | | 7.0 | | 15.0 | | | |
| 1,1,2-Trichloroethane | 0.3170<br>0.3219 | 0.3226 | 0.3388 | 0.3243 | 0.3317 | Ave | | 0.3260 | | | 0.1000 | 2.4 | | 15.0 | | | |
| Tetrachloroethene | 0.3819<br>0.3652 | 0.3715 | 0.3849 | 0.3699 | 0.3801 | Ave | | 0.3756 | | | 0.2000 | 2.1 | | 15.0 | | | |
| 1,3-Dichloropropane | 0.6750<br>0.6347 | 0.6523 | 0.6768 | 0.6516 | 0.6494 | Ave | | 0.6566 | | | | 2.5 | | 15.0 | | | |
| 2-Hexanone | 0.1779<br>0.1873 | 0.1707 | 0.1959 | 0.1832 | 0.1886 | Ave | | 0.1839 | | | 0.0500 | 4.8 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dibromochloromethane | 0.2426<br>0.3420 | 0.2434 | 0.2881 | 0.2960 | 0.3303 | Ave | | 0.2904 | | | | 14.4 | | 15.0 | | | |
| 1,2-Dibromoethane | 0.3215<br>0.3325 | 0.3183 | 0.3419 | 0.3251 | 0.3402 | Ave | | 0.3299 | | | | 3.0 | | 15.0 | | | |
| Chlorobenzene | 1.3869<br>1.2938 | 1.3415 | 1.3593 | 1.2905 | 1.3194 | Ave | | 1.3319 | | | 0.3000 | 2.8 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.3170<br>0.3986 | 0.3400 | 0.3683 | 0.3711 | 0.3999 | Ave | | 0.3658 | | | | 8.9 | | 15.0 | | | |
| Ethylbenzene | 0.7088<br>0.7002 | 0.6857 | 0.7184 | 0.6972 | 0.7117 | Ave | | 0.7037 | | | | 1.7 | | 15.0 | | | |
| m-Xylene & p-Xylene | 1.7254<br>1.7450 | 1.6919 | 1.7654 | 1.6850 | 1.7487 | Ave | | 1.7269 | | | | 1.9 | | 15.0 | | | |
| o-Xylene | 0.8614<br>0.8379 | 0.8285 | 0.8765 | 0.8245 | 0.8496 | Ave | | 0.8464 | | | | 2.4 | | 15.0 | | | |
| Styrene | 1.3641<br>1.5047 | 1.3699 | 1.4715 | 1.4375 | 1.5073 | Ave | | 1.4425 | | | 0.3000 | 4.4 | | 15.0 | | | |
| Bromoform | 0.1233<br>0.1962 | 0.1284 | 0.1486 | 0.1571 | 0.1803 | Lin1 | -0.109 | 0.1898 | | | 0.1000 | | | | 0.9940 | | 0.9900 |
| Isopropylbenzene | 2.0044<br>2.1719 | 2.0385 | 2.1403 | 2.0691 | 2.2120 | Ave | | 2.1060 | | | 0.1000 | 3.9 | | 15.0 | | | |
| Bromobenzene | 1.0386<br>0.9473 | 0.9586 | 0.9957 | 0.9686 | 0.9791 | Ave | | 0.9813 | | | | 3.3 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 0.7339<br>0.7541 | 0.7238 | 0.7477 | 0.7567 | 0.7581 | Ave | | 0.7457 | | | 0.3000 | 1.9 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.2416<br>0.2286 | 0.2178 | 0.2439 | 0.2276 | 0.2353 | Ave | | 0.2324 | | | | 4.2 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | 0.2160<br>0.2298 | 0.2002 | 0.2102 | 0.2187 | 0.2284 | Ave | | 0.2172 | | | | 5.2 | | 15.0 | | | |
| N-Propylbenzene | 1.0573<br>1.0569 | 1.0321 | 1.0773 | 1.0862 | 1.0836 | Ave | | 1.0656 | | | | 1.9 | | 15.0 | | | |
| 2-Chlorotoluene | 0.9709<br>0.9190 | 0.9662 | 0.9525 | 0.9301 | 0.9423 | Ave | | 0.9469 | | | | 2.1 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 3.2703<br>3.3967 | 3.1961 | 3.3564 | 3.2932 | 3.4464 | Ave | | 3.3265 | | | | 2.7 | | 15.0 | | | |
| 4-Chlorotoluene | 3.4379<br>3.3777 | 3.3433 | 3.4587 | 3.3511 | 3.4028 | Ave | | 3.3953 | | | | 1.4 | | 15.0 | | | |
| tert-Butylbenzene | 2.7549<br>2.7004 | 2.5816 | 2.7043 | 2.6702 | 2.7520 | Ave | | 2.6939 | | | | 2.4 | | 15.0 | | | |
| 1,2,4-Trimethylbenzene | 3.3781<br>3.5531 | 3.3438 | 3.4852 | 3.4900 | 3.5801 | Ave | | 3.4717 | | | | 2.7 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| sec-Butylbenzene | 3.6405<br>3.8669 | 3.6171 | 3.8143 | 3.8169 | 4.0039 | Ave | | 3.7933 | | | | 3.8 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.9907<br>1.8221 | 1.8786 | 1.9008 | 1.8807 | 1.8774 | Ave | | 1.8917 | | | 0.6000 | 2.9 | | 15.0 | | | |
| p-Isopropyltoluene | 3.1718<br>3.3680 | 3.1348 | 3.2894 | 3.3037 | 3.4483 | Ave | | 3.2860 | | | | 3.6 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 2.0885<br>1.8827 | 1.9679 | 1.9804 | 1.9148 | 1.9083 | Ave | | 1.9571 | | | 0.5000 | 3.8 | | 15.0 | | | |
| n-Butylbenzene | 2.5560<br>2.7774 | 2.6205 | 2.7472 | 2.7930 | 2.8960 | Ave | | 2.7317 | | | | 4.5 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.9442<br>1.7210 | 1.7936 | 1.8264 | 1.7788 | 1.7690 | Ave | | 1.8055 | | | 0.4000 | 4.2 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | 0.0900<br>0.1135 | 0.0828 | 0.0968 | 0.0997 | 0.1082 | Ave | | 0.0985 | | | 0.0500 | 11.5 | | 15.0 | | | |
| 1,2,4-Trichlorobenzene | 1.2023<br>1.0107 | 1.2116 | 1.1733 | 1.1572 | 1.1207 | Ave | | 1.1460 | | | 0.2000 | 6.4 | | 15.0 | | | |
| Hexachlorobutadiene | 0.5774<br>0.4251 | 0.5527 | 0.5232 | 0.5261 | 0.4932 | Ave | | 0.5163 | | | | 10.3 | | 15.0 | | | |
| Naphthalene | 2.2848<br>2.2596 | 2.3270 | 2.4025 | 2.4111 | 2.4450 | Ave | | 2.3550 | | | | 3.2 | | 15.0 | | | |
| 1,2,3-Trichlorobenzene | 1.0954<br>0.8485 | 1.0899 | 1.0634 | 1.0435 | 0.9857 | Ave | | 1.0211 | | | | 9.1 | | 15.0 | | | |
| Dibromofluoromethane (Surr) | 0.2849<br>0.2819 | 0.2702 | 0.3015 | 0.2862 | 0.2872 | Ave | | 0.2853 | | | | 3.5 | | 15.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | 0.3792<br>0.3401 | 0.3458 | 0.3615 | 0.3389 | 0.3376 | Ave | | 0.3505 | | | | 4.7 | | 15.0 | | | |
| Toluene-d8 (Surr) | 1.6874<br>1.6583 | 1.6036 | 1.7494 | 1.6376 | 1.6174 | Ave | | 1.6589 | | | | 3.2 | | 15.0 | | | |
| 4-Bromofluorobenzene (Surr) | 0.6307<br>0.6074 | 0.5837 | 0.6484 | 0.6004 | 0.6016 | Ave | | 0.6121 | | | | 3.8 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-210738/7 | UXC5227.D |
| Level 2 | STD8260 240-210738/6 | UXC5226.D |
| Level 3 | STD8260 240-210738/5 | UXC5225.D |
| Level 4 | STD8260 240-210738/4 | UXC5224.D |
| Level 5 | STD8260 240-210738/3 | UXC5223.D |
| Level 6 | STD8260 240-210738/2 | UXC5222A.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Dichlorodifluoromethane | FB | Ave | 69994<br>2968568 | 139619 | 372138 | 673816 | 1518235 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Chloromethane | FB | Ave | 87748<br>3083365 | 170524 | 420783 | 821450 | 1631821 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Vinyl chloride | FB | Ave | 71258<br>2881820 | 146004 | 378621 | 693671 | 1499386 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Butadiene | FB | Ave | 67249<br>2714311 | 137641 | 354913 | 644267 | 1414930 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Bromomethane | FB | Ave | 25554<br>929177 | 46647 | 142752 | 191122 | 549076 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Chloroethane | FB | Ave | 41750<br>1437795 | 78868 | 210854 | 366471 | 786655 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Dichlorofluoromethane | FB | Ave | 90041<br>3258061 | 164085 | 457144 | 778990 | 1782422 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Trichlorofluoromethane | FB | Ave | 58237<br>2512650 | 115835 | 319890 | 550869 | 1313251 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl ether | FB | Ave | 55376<br>1688348 | 91270 | 231842 | 475475 | 892922 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Acrolein | FB | Ave | 32432<br>1389519 | 61612 | 175522 | 346243 | 682062 | 5.00<br>200 | 10.0 | 25.0 | 50.0 | 100 |
| 1,1-Dichloroethene | FB | Ave | 55501<br>2353597 | 110800 | 286815 | 585373 | 1222106 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | FB | Ave | 38306<br>1837150 | 79207 | 208746 | 446581 | 968562 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Acetone | FB | Lin1 | 46991<br>1036072 | 69190 | 154585 | 274955 | 509043 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Iodomethane | FB | Ave | 96208<br>3852114 | 184743 | 478924 | 952532 | 1968866 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Carbon disulfide | FB | Ave | 165657<br>8105021 | 341907 | 907414 | 1888330 | 4067580 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 3-Chloro-1-propene | FB | Ave | 34429<br>1528096 | 66082 | 175712 | 374742 | 777919 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Methyl acetate | FB | Ave | 161589<br>6351704 | 316436 | 788908 | 1534412 | 3196421 | 5.00<br>200 | 10.0 | 25.0 | 50.0 | 100 |
| Methylene Chloride | FB | Lin1 | 169933<br>2604976 | 230171 | 441782 | 758268 | 1404532 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| tert-Butyl alcohol (TBA) | FB | Ave | 28243<br>1060473 | 55104 | 137053 | 277231 | 582101 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Methyl tert-butyl ether | FB | Ave | 140018<br>5722689 | 265059 | 697001 | 1391542 | 2899790 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| trans-1,2-Dichloroethene | FB | Ave | 62394<br>2494394 | 122598 | 314581 | 626109 | 1301937 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Acrylonitrile | FB | Ave | 168775<br>6989266 | 327470 | 847619 | 1682625 | 3513708 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Hexane | FB | Ave | 13860<br>617331 | 27863 | 74772 | 158931 | 331867 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1-Dichloroethane | FB | Ave | 114785<br>4624858 | 220177 | 587273 | 1149469 | 2378718 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Vinyl acetate | FB | Ave | 80620<br>3714612 | 155399 | 425331 | 865416 | 1846447 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2,2-Dichloropropane | FB | Ave | 58868<br>2563686 | 123228 | 323914 | 660593 | 1383113 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| cis-1,2-Dichloroethene | FB | Ave | 67276<br>2642242 | 129424 | 340423 | 669253 | 1367907 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Butanone (MEK) | FB | Ave | 38119<br>1446980 | 66526 | 181446 | 349299 | 718251 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Bromochloromethane | FB | Ave | 27982<br>1186680 | 57792 | 146678 | 297435 | 604822 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Tetrahydrofuran | FB | Lin1 | 31879<br>942339 | 55792 | 125662 | 233519 | 480109 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Chloroform | FB | Ave | 105796<br>4231988 | 200991 | 529540 | 1049547 | 2155600 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1,1-Trichloroethane | FB | Ave | 65867<br>3170356 | 144453 | 384623 | 781318 | 1684755 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Cyclohexane | FB | Ave | 102234<br>4585898 | 197683 | 537348 | 1140281 | 2405908 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1-Dichloropropene | FB | Ave | 78758<br>3396858 | 158611 | 419527 | 841750 | 1767419 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Carbon tetrachloride | FB | Ave | 54625<br>2830938 | 110857 | 303265 | 650617 | 1433806 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Isobutyl alcohol | FB | Ave | 28241<br>1066775 | 56556 | 134534 | 262630 | 562320 | 25.0<br>1000 | 50.0 | 125 | 250 | 500 |
| Benzene | FB | Ave | 277110<br>10921918 | 518188 | 1342060 | 2637881 | 5557555 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| 1,2-Dichloroethane | FB | Ave | 79153<br>3100592 | 152324 | 397554 | 776612 | 1600630 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| n-Heptane | FB | Ave | 14769<br>572892 | 24035 | 65355 | 145300 | 310101 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Trichloroethene | FB | Ave | 62443<br>2555067 | 121983 | 322045 | 636127 | 1325685 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Methylcyclohexane | FB | Ave | 80182<br>3872636 | 169040 | 449695 | 965944 | 2071671 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dichloropropane | FB | Ave | 59802<br>2528018 | 121453 | 310531 | 624994 | 1278043 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Dibromomethane | FB | Ave | 30272<br>1229909 | 56615 | 148707 | 295275 | 632539 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,4-Dioxane | FB | Ave | 9085<br>346500 | 17434 | 43294 | 90456 | 196830 | 20.0<br>800 | 40.0 | 100 | 200 | 400 |
| Bromodichloromethane | FB | Ave | 56307<br>2874987 | 109399 | 312560 | 656783 | 1425790 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Chloroethyl vinyl ether | FB | Ave | 56171<br>2577398 | 111631 | 303102 | 619938 | 1305368 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| cis-1,3-Dichloropropene | FB | Ave | 67826<br>3696869 | 143257 | 398000 | 835323 | 1827980 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | FB | Ave | 70676<br>2999002 | 129333 | 341762 | 704866 | 1477206 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Toluene | CBZ | Ave | 260892<br>10830971 | 511399 | 1302798 | 2586870 | 5499121 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| trans-1,3-Dichloropropene | CBZ | Ave | 48815<br>2879764 | 104714 | 301173 | 629964 | 1413295 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl methacrylate | CBZ | Ave | 46794<br>2397003 | 96464 | 267551 | 561135 | 1193908 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1,2-Trichloroethane | CBZ | Ave | 39523<br>1734782 | 81281 | 216506 | 429312 | 898262 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Tetrachloroethene | CBZ | Ave | 47609<br>1968241 | 93612 | 245987 | 489679 | 1029269 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,3-Dichloropropane | CBZ | Ave | 84161<br>3421307 | 164354 | 432575 | 862483 | 1758568 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Hexanone | CBZ | Ave | 44351<br>2018747 | 86032 | 250356 | 485141 | 1021566 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Dibromochloromethane | CBZ | Ave | 30248<br>1843569 | 61336 | 184148 | 391874 | 894516 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dibromoethane | CBZ | Ave | 40077<br>1791978 | 80199 | 218492 | 430310 | 921303 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Chlorobenzene | CBZ | Ave | 172909<br>6973434 | 337996 | 868719 | 1708229 | 3573219 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| 1,1,1,2-Tetrachloroethane | CBZ | Ave | 39522<br>2148347 | 85665 | 235415 | 491177 | 1083094 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethylbenzene | CBZ | Ave | 88364<br>3774329 | 172765 | 459128 | 922946 | 1927335 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| m-Xylene & p-Xylene | CBZ | Ave | 215117<br>9405749 | 426271 | 1128260 | 2230521 | 4735552 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| o-Xylene | CBZ | Ave | 107394<br>4516212 | 208753 | 560164 | 1091371 | 2300904 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Styrene | CBZ | Ave | 170064<br>8110542 | 345149 | 940436 | 1902845 | 4082071 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Bromoform | CBZ | Lin1 | 15373<br>1057395 | 32348 | 95000 | 207939 | 488341 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Isopropylbenzene | CBZ | Ave | 249899<br>11706432 | 513612 | 1367862 | 2738887 | 5990259 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Bromobenzene | DCB | Ave | 67655<br>2699119 | 126334 | 341138 | 659380 | 1410057 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,1,2,2-Tetrachloroethane | DCB | Ave | 47806<br>2148675 | 95393 | 256175 | 515138 | 1091773 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2,3-Trichloropropane | DCB | Ave | 15740<br>651190 | 28698 | 83557 | 154970 | 338794 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| trans-1,4-Dichloro-2-butene | DCB | Ave | 14068<br>654679 | 26383 | 72027 | 148910 | 328971 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| N-Propylbenzene | DCB | Ave | 68871<br>3011323 | 136029 | 369098 | 739451 | 1560563 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Chlorotoluene | DCB | Ave | 63247<br>2618465 | 127341 | 326359 | 633207 | 1357052 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,3,5-Trimethylbenzene | DCB | Ave | 213034<br>9677816 | 421217 | 1150007 | 2241918 | 4963247 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 4-Chlorotoluene | DCB | Ave | 223946<br>9623665 | 440627 | 1185064 | 2281347 | 4900450 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| tert-Butylbenzene | DCB | Ave | 179458<br>7693891 | 340237 | 926572 | 1817797 | 3963274 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2,4-Trimethylbenzene | DCB | Ave | 220056<br>10123377 | 440692 | 1194129 | 2375931 | 5155790 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| sec-Butylbenzene | DCB | Ave | 237149<br>11017467 | 476701 | 1306888 | 2598481 | 5766100 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,3-Dichlorobenzene | DCB | Ave | 129674<br>5191608 | 247588 | 651276 | 1280328 | 2703641 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| p-Isopropyltoluene | DCB | Ave | 206615<br>9596155 | 413139 | 1127050 | 2249065 | 4965916 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,4-Dichlorobenzene | DCB | Ave | 136050<br>5364042 | 259349 | 678527 | 1303553 | 2748121 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 13:08 Calibration End Date: 12/14/2015 15:01 Calibration ID: 32097

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| n-Butylbenzene | DCB | Ave | 166504<br>7913388 | 345364 | 941255 | 1901423 | 4170653 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dichlorobenzene | DCB | Ave | 126645<br>4903421 | 236378 | 625774 | 1210963 | 2547569 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dibromo-3-Chloropropane | DCB | Ave | 5861<br>323252 | 10913 | 33170 | 67842 | 155810 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2,4-Trichlorobenzene | DCB | Ave | 78318<br>2879725 | 159681 | 402011 | 787791 | 1613918 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Hexachlorobutadiene | DCB | Ave | 37612<br>1211136 | 72848 | 179268 | 358150 | 710309 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Naphthalene | DCB | Ave | 148836<br>6438081 | 306684 | 823174 | 1641414 | 3521077 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2,3-Trichlorobenzene | DCB | Ave | 71354<br>2417629 | 143641 | 364350 | 710386 | 1419594 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Dibromofluoromethane (Surr) | FB | Ave | 47893<br>2052047 | 92758 | 263782 | 520230 | 1045565 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | FB | Ave | 63759<br>2475424 | 118695 | 316270 | 615898 | 1229201 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Toluene-d8 (Surr) | CBZ | Ave | 210371<br>8938007 | 404025 | 1118072 | 2167677 | 4380013 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 4-Bromofluorobenzene (Surr) | CBZ | Ave | 78632<br>3273800 | 147068 | 414424 | 794830 | 1629282 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin1 = Linear 1/conc ISTD |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 15:47 Calibration End Date: 12/14/2015 17:40 Calibration ID: 32099

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD 240-210738/13 | UXC5234.D |
| Level 2 | STD 240-210738/12 | UXC5233.D |
| Level 3 | STD 240-210738/11 | UXC5232.D |
| Level 4 | STD 240-210738/10 | UXC5231.D |
| Level 5 | STD 240-210738/9 | UXC5230.D |
| Level 6 | STD 240-210738/8 | UXC5229.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Acetonitrile | 0.0365<br>0.0317 | 0.0300 | 0.0294 | 0.0290 | 0.0359 | Ave | | 0.0321 | | | | 10.4 | | 15.0 | | | |
| Diisopropyl ether | 0.3037<br>0.3132 | 0.2709 | 0.2911 | 0.2858 | 0.3539 | Ave | | 0.3031 | | | | 9.5 | | 15.0 | | | |
| 2-Chloro-1,3-butadiene | 0.5766<br>0.5821 | 0.4936 | 0.5222 | 0.5190 | 0.6414 | Ave | | 0.5558 | | | | 9.8 | | 15.0 | | | |
| Ethyl t-butyl ether | 0.8978<br>0.9605 | 0.7949 | 0.8552 | 0.8432 | 1.0631 | Ave | | 0.9024 | | | | 10.7 | | 15.0 | | | |
| Ethyl acetate | 0.2184<br>0.2097 | 0.1742 | 0.1936 | 0.1876 | 0.2344 | Ave | | 0.2030 | | | | 10.8 | | 15.0 | | | |
| Propionitrile | 0.0343<br>0.0353 | 0.0303 | 0.0330 | 0.0324 | 0.0400 | Ave | | 0.0342 | | | | 9.7 | | 15.0 | | | |
| Methacrylonitrile | 0.1640<br>0.1627 | 0.1420 | 0.1484 | 0.1464 | 0.1822 | Ave | | 0.1576 | | | | 9.5 | | 15.0 | | | |
| Tert-amyl methyl ether | 0.7246<br>0.7757 | 0.6354 | 0.6812 | 0.6816 | 0.8534 | Ave | | 0.7253 | | | | 10.8 | | 15.0 | | | |
| n-Butanol | 0.0044<br>0.0048 | 0.0040 | 0.0043 | 0.0042 | 0.0056 | Ave | | 0.0046 | | | | 13.1 | | 15.0 | | | |
| Ethyl acrylate | 0.2503<br>0.2916 | 0.2326 | 0.2532 | 0.2494 | 0.3195 | Ave | | 0.2661 | | | | 12.3 | | 15.0 | | | |
| Methyl methacrylate | 0.1924<br>0.2149 | 0.1720 | 0.1912 | 0.1885 | 0.2390 | Ave | | 0.1997 | | | | 11.8 | | 15.0 | | | |
| 2-Nitropropane | 0.0253<br>0.0344 | 0.0215 | 0.0251 | 0.0259 | 0.0357 | Qua | -0.070 | 0.0349 | 0.0000078 | | | | | | 0.9960 | | 0.9900 |
| n-Butyl acetate | 0.2685<br>0.3405 | 0.2390 | 0.2761 | 0.2795 | 0.3735 | Lin1 | -0.158 | 0.3457 | | | | | | | 0.9910 | | 0.9900 |
| 1-Chlorohexane | 0.5485<br>0.5838 | 0.4788 | 0.5134 | 0.5231 | 0.6338 | Ave | | 0.5469 | | | | 10.1 | | 15.0 | | | |
| Cyclohexanone | 0.0158<br>0.0171 | 0.0149 | 0.0156 | 0.0157 | 0.0200 | Ave | | 0.0165 | | | | 11.3 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 15:47 Calibration End Date: 12/14/2015 17:40 Calibration ID: 32099

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,2,3-Trimethylbenzene | 3.5383<br>3.8182 | 3.0588 | 3.3758 | 3.3699 | 4.2665 | Ave | | 3.5712 | | | | 11.8 | | 15.0 | | | |
| 1,3,5-Trichlorobenzene | 1.2998<br>1.2091 | 1.1604 | 1.1860 | 1.1935 | 1.4476 | Ave | | 1.2494 | | | | 8.7 | | 15.0 | | | |
| 2-Methylnaphthalene | 1.4202<br>+++++ | 1.4607 | 1.3806 | 1.4983 | 1.6153 | Ave | | 1.4750 | | | | 6.1 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 15:47 Calibration End Date: 12/14/2015 17:40 Calibration ID: 32099

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD 240-210738/13 | UXC5234.D |
| Level 2 | STD 240-210738/12 | UXC5233.D |
| Level 3 | STD 240-210738/11 | UXC5232.D |
| Level 4 | STD 240-210738/10 | UXC5231.D |
| Level 5 | STD 240-210738/9 | UXC5230.D |
| Level 6 | STD 240-210738/8 | UXC5229.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Acetonitrile | FB | Ave | 58710<br>2161234 | 107618 | 269440 | 530427 | 1089443 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Diisopropyl ether | FB | Ave | 48874<br>2135571 | 97225 | 266599 | 523206 | 1073436 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Chloro-1,3-butadiene | FB | Ave | 92811<br>3968725 | 177164 | 478135 | 950073 | 1945304 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl t-butyl ether | FB | Ave | 144495<br>6549049 | 285292 | 783117 | 1543592 | 3224142 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Ethyl acetate | FB | Ave | 70314<br>2859248 | 125066 | 354507 | 687008 | 1421812 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| Propionitrile | FB | Ave | 55145<br>2403671 | 108661 | 302198 | 593215 | 1214198 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Methacrylonitrile | FB | Ave | 263884<br>11093173 | 509496 | 1359229 | 2679197 | 5527214 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| Tert-amyl methyl ether | FB | Ave | 116618<br>5288533 | 228068 | 623735 | 1247683 | 2588171 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| n-Butanol | FB | Ave | 17662<br>824245 | 36094 | 97748 | 191701 | 427606 | 25.0<br>1000 | 50.0 | 125 | 250 | 500 |
| Ethyl acrylate | FB | Ave | 40285<br>1988241 | 83488 | 231812 | 456560 | 969022 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Methyl methacrylate | FB | Ave | 61926<br>2930581 | 123470 | 350113 | 690199 | 1449638 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| 2-Nitropropane | FB | Qua | 8137<br>469745 | 15450 | 45918 | 94990 | 216285 | 2.00<br>80.0 | 4.00 | 10.0 | 20.0 | 40.0 |
| n-Butyl acetate | FB | Lin1 | 43214<br>2321703 | 85795 | 252849 | 511700 | 1132859 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 1-Chlorohexane | CBZ | Ave | 65035<br>2917244 | 125109 | 340397 | 691011 | 1421577 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| Cyclohexanone | DCB | Ave | 9422<br>445155 | 19855 | 52906 | 107609 | 231275 | 10.0<br>400 | 20.0 | 50.0 | 100 | 200 |
| 1,2,3-Trimethylbenzene | DCB | Ave | 211554<br>9941527 | 408349 | 1143850 | 2303984 | 4923338 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210738

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/14/2015 15:47 Calibration End Date: 12/14/2015 17:40 Calibration ID: 32099

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| 1,3,5-Trichlorobenzene | DCB | Ave | 77714<br>3148084 | 154908 | 401857 | 816024 | 1670425 | 1.00<br>40.0 | 2.00 | 5.00 | 10.0 | 20.0 |
| 2-Methylnaphthalene | DCB | Ave | 169829<br>+++++ | 389994 | 935608 | 2048722 | 3728051 | 2.00<br>+++++ | 4.00 | 10.0 | 20.0 | 40.0 |

Curve Type Legend:

Ave = Average ISTD
Lin1 = Linear 1/conc ISTD
Qua = Quadratic ISTD

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-212606/21 Calibration Date: 12/30/2015 18:56

Instrument ID: A3UX14 Calib Start Date: 12/30/2015 12:29

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 12/30/2015 15:21

Lab File ID: 147318.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.3057 | 0.3115 | 0.1000 | 0.0204 | 0.0200 | 1.9 | 20.0 |
| Chloromethane | Ave | 0.4683 | 0.4789 | 0.1000 | 0.0205 | 0.0200 | 2.3 | 20.0 |
| Vinyl chloride | Ave | 0.3108 | 0.3139 | 0.1000 | 0.0202 | 0.0200 | 1.0 | 20.0 |
| Butadiene | Ave | 0.3624 | 0.3725 | | 0.0206 | 0.0200 | 2.8 | 20.0 |
| Bromomethane | Ave | 0.1379 | 0.1264 | 0.0500 | 0.0183 | 0.0200 | -8.3 | 20.0 |
| Chloroethane | Ave | 0.1675 | 0.1628 | 0.0500 | 0.0194 | 0.0200 | -2.8 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3382 | 0.3132 | | 0.0185 | 0.0200 | -7.4 | 20.0 |
| Trichlorofluoromethane | Ave | 0.3441 | 0.3707 | 0.1000 | 0.0215 | 0.0200 | 7.7 | 20.0 |
| Ethyl ether | Ave | 0.2190 | 0.2207 | | 0.0504 | 0.0500 | 0.8 | 20.0 |
| Acrolein | Lin1 | | 0.0305 | | 0.229 | 0.250 | -8.4 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2482 | 0.2779 | 0.1000 | 0.0560 | 0.0500 | 12.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2084 | 0.2215 | 0.0500 | 0.0532 | 0.0500 | 6.3 | 20.0 |
| Acetone | Qua | | 0.0773 | 0.0500 | 0.0837 | 0.100 | -16.3 | 20.0 |
| Iodomethane | Ave | 0.3700 | 0.4022 | | 0.0543 | 0.0500 | 8.7 | 20.0 |
| Carbon disulfide | Ave | 0.7420 | 0.7935 | 0.1000 | 0.0535 | 0.0500 | 6.9 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1467 | 0.1628 | | 0.0555 | 0.0500 | 11.0 | 20.0 |
| Methyl acetate | Ave | 0.1860 | 0.1851 | 0.1000 | 0.249 | 0.250 | -0.5 | 20.0 |
| Methylene Chloride | Qua | | 0.2983 | 0.1000 | 0.0523 | 0.0500 | 4.7 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0247 | 0.0223 | | 0.451 | 0.500 | -9.7 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2941 | 0.3151 | 0.1000 | 0.0536 | 0.0500 | 7.1 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.6281 | 0.6378 | 0.1000 | 0.0508 | 0.0500 | 1.5 | 20.0 |
| Acrylonitrile | Ave | 0.1038 | 0.1028 | | 0.495 | 0.500 | -0.9 | 20.0 |
| Hexane | Ave | 0.0811 | 0.0885 | | 0.0546 | 0.0500 | 9.2 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.6440 | 0.6644 | 0.1000 | 0.0516 | 0.0500 | 3.2 | 20.0 |
| Vinyl acetate | Ave | 0.0331 | 0.0383 | | 0.0577 | 0.0500 | 15.5 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2876 | 0.2907 | | 0.0505 | 0.0500 | 1.1 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3014 | 0.3238 | 0.1000 | 0.0537 | 0.0500 | 7.4 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.1130 | 0.1044 | 0.0500 | 0.0924 | 0.100 | -7.6 | 20.0 |
| Bromochloromethane | Ave | 0.1197 | 0.1226 | | 0.0512 | 0.0500 | 2.5 | 20.0 |
| Tetrahydrofuran | Ave | 0.0755 | 0.0706 | | 0.0935 | 0.100 | -6.5 | 20.0 |
| Chloroform | Ave | 0.4940 | 0.5120 | 0.2000 | 0.0518 | 0.0500 | 3.6 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.4136 | 0.4561 | 0.1000 | 0.0551 | 0.0500 | 10.3 | 20.0 |
| Cyclohexane | Ave | 0.7887 | 0.8710 | 0.1000 | 0.0552 | 0.0500 | 10.4 | 20.0 |
| Carbon tetrachloride | Ave | 0.3312 | 0.3888 | 0.1000 | 0.0587 | 0.0500 | 17.4 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3643 | 0.4135 | | 0.0568 | 0.0500 | 13.5 | 20.0 |
| Benzene | Ave | 1.106 | 1.162 | 0.5000 | 0.0525 | 0.0500 | 5.1 | 20.0 |
| Isobutyl alcohol | Ave | 0.0081 | 0.0077 | | 1.19 | 1.25 | -4.8 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.4002 | 0.4237 | 0.1000 | 0.0529 | 0.0500 | 5.9 | 20.0 |
| n-Heptane | Ave | 0.0755 | 0.0837 | | 0.0554 | 0.0500 | 10.9 | 20.0 |
| Trichloroethene | Ave | 0.2754 | 0.2984 | 0.2000 | 0.0542 | 0.0500 | 8.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-212606/21 Calibration Date: 12/30/2015 18:56

Instrument ID: A3UX14 Calib Start Date: 12/30/2015 12:29

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 12/30/2015 15:21

Lab File ID: 147318.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.5152 | 0.5652 | 0.1000 | 0.0549 | 0.0500 | 9.7 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.3148 | 0.3351 | 0.1000 | 0.0532 | 0.0500 | 6.4 | 20.0 |
| Dibromomethane | Ave | 0.1235 | 0.1328 | | 0.0538 | 0.0500 | 7.6 | 20.0 |
| 1,4-Dioxane | Ave | 0.0019 | 0.0018 | | 0.979 | 1.00 | -2.1 | 20.0 |
| Bromodichloromethane | Ave | 0.2898 | 0.3175 | 0.1000 | 0.0548 | 0.0500 | 9.6 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0951 | 0.0998 | | 0.0525 | 0.0500 | 5.0 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3353 | 0.3613 | 0.1500 | 0.0539 | 0.0500 | 7.8 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3685 | 0.3631 | 0.0500 | 0.0985 | 0.100 | -1.5 | 20.0 |
| Toluene | Ave | 1.863 | 1.943 | 0.4000 | 0.0521 | 0.0500 | 4.3 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.4211 | 0.4412 | 0.1000 | 0.0524 | 0.0500 | 4.8 | 20.0 |
| Ethyl methacrylate | Ave | 0.3247 | 0.3528 | | 0.0543 | 0.0500 | 8.7 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2910 | 0.2824 | 0.1000 | 0.0485 | 0.0500 | -2.9 | 20.0 |
| Tetrachloroethene | Ave | 0.3443 | 0.3443 | 0.2000 | 0.0500 | 0.0500 | 0.0 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4900 | 0.4778 | | 0.0488 | 0.0500 | -2.5 | 20.0 |
| 2-Hexanone | Ave | 0.2523 | 0.2214 | 0.0500 | 0.0877 | 0.100 | -12.3 | 20.0 |
| Dibromochloromethane | Ave | 0.2828 | 0.2872 | | 0.0508 | 0.0500 | 1.5 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2561 | 0.2669 | | 0.0521 | 0.0500 | 4.2 | 20.0 |
| Chlorobenzene | Ave | 1.077 | 1.085 | 0.3000 | 0.0504 | 0.0500 | 0.8 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3503 | 0.3843 | | 0.0548 | 0.0500 | 9.7 | 20.0 |
| Ethylbenzene | Ave | 0.6092 | 0.6183 | | 0.0507 | 0.0500 | 1.5 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.7344 | 0.7523 | | 0.0512 | 0.0500 | 2.4 | 20.0 |
| o-Xylene | Ave | 0.7252 | 0.7613 | | 0.0525 | 0.0500 | 5.0 | 20.0 |
| Styrene | Ave | 1.080 | 1.120 | 0.3000 | 0.0519 | 0.0500 | 3.7 | 20.0 |
| Bromoform | Ave | 0.1467 | 0.1520 | 0.1000 | 0.0518 | 0.0500 | 3.6 | 20.0 |
| Isopropylbenzene | Ave | 2.047 | 2.179 | 0.1000 | 0.0532 | 0.0500 | 6.5 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.6475 | 0.6808 | 0.3000 | 0.0526 | 0.0500 | 5.1 | 20.0 |
| Bromobenzene | Ave | 0.8490 | 0.8832 | | 0.0520 | 0.0500 | 4.0 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1864 | 0.2058 | | 0.0552 | 0.0500 | 10.4 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2792 | 0.2814 | | 0.0504 | 0.0500 | 0.8 | 20.0 |
| N-Propylbenzene | Ave | 1.038 | 1.134 | | 0.0546 | 0.0500 | 9.2 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8878 | 0.9391 | | 0.0529 | 0.0500 | 5.8 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.363 | 3.721 | | 0.0553 | 0.0500 | 10.6 | 20.0 |
| 4-Chlorotoluene | Ave | 0.8863 | 0.9031 | | 0.0509 | 0.0500 | 1.9 | 20.0 |
| tert-Butylbenzene | Ave | 3.023 | 3.386 | | 0.0560 | 0.0500 | 12.0 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.372 | 3.634 | | 0.0539 | 0.0500 | 7.8 | 20.0 |
| sec-Butylbenzene | Ave | 4.417 | 4.872 | | 0.0552 | 0.0500 | 10.3 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.629 | 1.607 | 0.6000 | 0.0493 | 0.0500 | -1.3 | 20.0 |
| p-Isopropyltoluene | Ave | 3.661 | 4.030 | | 0.0550 | 0.0500 | 10.1 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.635 | 1.579 | 0.5000 | 0.0483 | 0.0500 | -3.5 | 20.0 |
| n-Butylbenzene | Ave | 3.550 | 3.760 | | 0.0530 | 0.0500 | 5.9 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.591 | 1.546 | 0.4000 | 0.0486 | 0.0500 | -2.8 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-212606/21  Calibration Date: 12/30/2015 18:56

Instrument ID: A3UX14  Calib Start Date: 12/30/2015 12:29

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 12/30/2015 15:21

Lab File ID: 147318.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.1066 | 0.1091 | 0.0500 | 0.0512 | 0.0500 | 2.3 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 1.258 | 1.170 | 0.2000 | 0.0465 | 0.0500 | -7.0 | 20.0 |
| Hexachlorobutadiene | Ave | 0.8161 | 0.8312 | | 0.0509 | 0.0500 | 1.8 | 20.0 |
| Naphthalene | Ave | 2.465 | 2.323 | | 0.0471 | 0.0500 | -5.8 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 1.197 | 1.075 | | 0.0449 | 0.0500 | -10.2 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2305 | 0.2194 | | 0.0476 | 0.0500 | -4.8 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.3100 | 0.2860 | | 0.0461 | 0.0500 | -7.7 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.607 | 1.491 | | 0.0464 | 0.0500 | -7.2 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.139 | 1.061 | | 0.0466 | 0.0500 | -6.8 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-212606/22  Calibration Date: 12/30/2015 19:17

Instrument ID: A3UX14  Calib Start Date: 12/30/2015 12:29

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 12/30/2015 15:21

Lab File ID: 147319.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dibromofluoromethane (Surr) | Ave | 0.2305 | 0.2092 | | 0.0454 | 0.0500 | -9.3 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.3100 | 0.2793 | | 0.0451 | 0.0500 | -9.9 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.607 | 1.551 | | 0.0482 | 0.0500 | -3.5 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.139 | 0.998 | | 0.0438 | 0.0500 | -12.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-212606/22  Calibration Date: 12/30/2015 19:17

Instrument ID: A3UX14  Calib Start Date: 12/30/2015 15:42

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 12/30/2015 18:34

Lab File ID: 147319.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0319 | 0.0310 | | 0.485 | 0.500 | -3.1 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.6633 | 0.7274 | | 0.0548 | 0.0500 | 9.7 | 20.0 |
| Diisopropyl ether | Ave | 0.2233 | 0.2521 | | 0.0564 | 0.0500 | 12.9 | 20.0 |
| Ethyl t-butyl ether | Ave | 1.013 | 1.078 | | 0.0532 | 0.0500 | 6.4 | 20.0 |
| Ethyl acetate | Ave | 0.2088 | 0.2111 | | 0.101 | 0.100 | 1.1 | 20.0 |
| Propionitrile | Ave | 0.0384 | 0.0380 | | 0.494 | 0.500 | -1.2 | 20.0 |
| Methacrylonitrile | Ave | 0.1427 | 0.1516 | | 0.531 | 0.500 | 6.2 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.6388 | 0.6765 | | 0.0530 | 0.0500 | 5.9 | 20.0 |
| n-Butanol | Ave | 0.0118 | 0.0124 | | 1.32 | 1.25 | 5.4 | 20.0 |
| Methyl methacrylate | Ave | 0.1860 | 0.1937 | | 0.104 | 0.100 | 4.2 | 20.0 |
| 2-Nitropropane | Ave | 0.0380 | 0.0368 | | 0.0969 | 0.100 | -3.1 | 20.0 |
| 1-Chlorohexane | Ave | 0.5064 | 0.5283 | | 0.0522 | 0.0500 | 4.3 | 20.0 |
| Cyclohexanone | Ave | 0.0142 | 0.0148 | | 0.522 | 0.500 | 4.4 | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.279 | 3.616 | | 0.0551 | 0.0500 | 10.3 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.290 | 1.388 | | 0.0538 | 0.0500 | 7.6 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.461 | 1.513 | | 0.104 | 0.100 | 3.5 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-210738/14 Calibration Date: 12/14/2015 15:24

Instrument ID: A3UX15 Calib Start Date: 12/14/2015 13:08

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 12/14/2015 15:01

Lab File ID: UXC5228.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.4073 | 0.4509 | 0.1000 | 0.0111 | 0.0100 | 10.7 | 20.0 |
| Chloromethane | Ave | 0.4706 | 0.4781 | 0.1000 | 0.0102 | 0.0100 | 1.6 | 20.0 |
| Vinyl chloride | Ave | 0.4119 | 0.4351 | 0.1000 | 0.0106 | 0.0100 | 5.6 | 20.0 |
| Butadiene | Ave | 0.3871 | 0.4016 | | 0.0104 | 0.0100 | 3.7 | 20.0 |
| Bromomethane | Ave | 0.1391 | 0.1718 | 0.0500 | 0.0123 | 0.0100 | 23.5* | 20.0 |
| Chloroethane | Ave | 0.2224 | 0.2327 | 0.0500 | 0.0105 | 0.0100 | 4.6 | 20.0 |
| Dichlorofluoromethane | Ave | 0.4837 | 0.5307 | | 0.0110 | 0.0100 | 9.7 | 20.0 |
| Trichlorofluoromethane | Ave | 0.3431 | 0.3875 | 0.1000 | 0.0113 | 0.0100 | 12.9 | 20.0 |
| Ethyl ether | Ave | 0.2665 | 0.2619 | | 0.00983 | 0.0100 | -1.7 | 20.0 |
| Acrolein | Ave | 0.0381 | 0.0404 | | 0.0531 | 0.0500 | 6.3 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.3270 | 0.3244 | 0.1000 | 0.00992 | 0.0100 | -0.8 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2436 | 0.2656 | 0.0500 | 0.0109 | 0.0100 | 9.0 | 20.0 |
| Acetone | Lin1 | | 0.0778 | 0.0500 | 0.0206 | 0.0200 | 2.8 | 20.0 |
| Iodomethane | Ave | 0.5420 | 0.5628 | | 0.0104 | 0.0100 | 3.8 | 20.0 |
| Carbon disulfide | Ave | 1.048 | 1.058 | 0.1000 | 0.0101 | 0.0100 | 0.9 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.2047 | 0.1982 | | 0.00968 | 0.0100 | -3.2 | 20.0 |
| Methyl acetate | Ave | 0.1793 | 0.1854 | 0.1000 | 0.0517 | 0.0500 | 3.4 | 20.0 |
| Methylene Chloride | Lin1 | | 0.4220 | 0.1000 | 0.0102 | 0.0100 | 2.3 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0157 | 0.0151 | | 0.0960 | 0.100 | -4.0 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7917 | 0.8101 | 0.1000 | 0.0102 | 0.0100 | 2.3 | 20.0 |
| Acrylonitrile | Ave | 0.0963 | 0.0999 | | 0.104 | 0.100 | 3.7 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.3554 | 0.3687 | 0.1000 | 0.0104 | 0.0100 | 3.7 | 20.0 |
| Hexane | Ave | 0.0854 | 0.0919 | | 0.0108 | 0.0100 | 7.6 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.6528 | 0.6763 | 0.1000 | 0.0104 | 0.0100 | 3.6 | 20.0 |
| Vinyl acetate | Ave | 0.4853 | 0.4614 | | 0.00951 | 0.0100 | -4.9 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.3625 | 0.3701 | | 0.0102 | 0.0100 | 2.1 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3789 | 0.3915 | 0.1000 | 0.0103 | 0.0100 | 3.3 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.1014 | 0.1013 | 0.0500 | 0.0200 | 0.0200 | -0.0 | 20.0 |
| Bromochloromethane | Ave | 0.1659 | 0.1712 | | 0.0103 | 0.0100 | 3.2 | 20.0 |
| Tetrahydrofuran | Lin1 | | 0.0687 | | 0.0205 | 0.0200 | 2.7 | 20.0 |
| Chloroform | Ave | 0.5952 | 0.6124 | 0.2000 | 0.0103 | 0.0100 | 2.9 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.4301 | 0.4572 | 0.1000 | 0.0106 | 0.0100 | 6.3 | 20.0 |
| Cyclohexane | Ave | 0.6194 | 0.6704 | 0.1000 | 0.0108 | 0.0100 | 8.2 | 20.0 |
| Carbon tetrachloride | Ave | 0.3559 | 0.3858 | 0.1000 | 0.0108 | 0.0100 | 8.4 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.4709 | 0.5435 | | 0.0115 | 0.0100 | 15.4 | 20.0 |
| Isobutyl alcohol | Ave | 0.0062 | 0.0058 | | 0.233 | 0.250 | -6.8 | 20.0 |
| Benzene | Ave | 1.528 | 1.550 | 0.5000 | 0.0101 | 0.0100 | 1.4 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.4436 | 0.4534 | 0.1000 | 0.0102 | 0.0100 | 2.2 | 20.0 |
| n-Heptane | Ave | 0.0794 | 0.0855 | | 0.0108 | 0.0100 | 7.6 | 20.0 |
| Trichloroethene | Ave | 0.3600 | 0.3771 | 0.2000 | 0.0105 | 0.0100 | 4.7 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-210738/14 Calibration Date: 12/14/2015 15:24

Instrument ID: A3UX15 Calib Start Date: 12/14/2015 13:08

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 12/14/2015 15:01

Lab File ID: UXC5228.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.5193 | 0.5721 | 0.1000 | 0.0110 | 0.0100 | 10.2 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.3511 | 0.3683 | 0.1000 | 0.0105 | 0.0100 | 4.9 | 20.0 |
| Dibromomethane | Ave | 0.1700 | 0.1755 | | 0.0103 | 0.0100 | 3.2 | 20.0 |
| 1,4-Dioxane | Ave | 0.0025 | 0.0024 | | 0.189 | 0.200 | -5.6 | 20.0 |
| Bromodichloromethane | Ave | 0.3598 | 0.3784 | 0.1000 | 0.0105 | 0.0100 | 5.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.1716 | 0.1785 | | 0.0104 | 0.0100 | 4.0 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.4575 | 0.4934 | 0.1500 | 0.0108 | 0.0100 | 7.8 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.1994 | 0.2078 | 0.0500 | 0.0208 | 0.0200 | 4.2 | 20.0 |
| Toluene | Ave | 2.026 | 2.085 | 0.4000 | 0.0103 | 0.0100 | 2.9 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.4684 | 0.5450 | 0.1000 | 0.0116 | 0.0100 | 16.4 | 20.0 |
| Ethyl methacrylate | Ave | 0.4144 | 0.4593 | | 0.0111 | 0.0100 | 10.8 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.3260 | 0.3420 | 0.1000 | 0.0105 | 0.0100 | 4.9 | 20.0 |
| Tetrachloroethene | Ave | 0.3756 | 0.3897 | 0.2000 | 0.0104 | 0.0100 | 3.8 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.6566 | 0.6693 | | 0.0102 | 0.0100 | 1.9 | 20.0 |
| 2-Hexanone | Ave | 0.1839 | 0.1882 | 0.0500 | 0.0205 | 0.0200 | 2.3 | 20.0 |
| Dibromochloromethane | Ave | 0.2904 | 0.3073 | | 0.0106 | 0.0100 | 5.8 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.3299 | 0.3412 | | 0.0103 | 0.0100 | 3.4 | 20.0 |
| Chlorobenzene | Ave | 1.332 | 1.347 | 0.3000 | 0.0101 | 0.0100 | 1.1 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3658 | 0.3834 | | 0.0105 | 0.0100 | 4.8 | 20.0 |
| Ethylbenzene | Ave | 0.7037 | 0.7349 | | 0.0104 | 0.0100 | 4.4 | 20.0 |
| m-Xylene & p-Xylene | Ave | 1.727 | 1.770 | | 0.0102 | 0.0100 | 2.5 | 20.0 |
| o-Xylene | Ave | 0.8464 | 0.8791 | | 0.0104 | 0.0100 | 3.9 | 20.0 |
| Styrene | Ave | 1.443 | 1.445 | 0.3000 | 0.0100 | 0.0100 | 0.2 | 20.0 |
| Bromoform | Lin1 | | 0.1566 | 0.1000 | 0.00883 | 0.0100 | -11.7 | 20.0 |
| Isopropylbenzene | Ave | 2.106 | 2.201 | 0.1000 | 0.0104 | 0.0100 | 4.5 | 20.0 |
| Bromobenzene | Ave | 0.9813 | 1.002 | | 0.0102 | 0.0100 | 2.1 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.7457 | 0.7742 | 0.3000 | 0.0104 | 0.0100 | 3.8 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.2324 | 0.2558 | | 0.0110 | 0.0100 | 10.1 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2172 | 0.2383 | | 0.0110 | 0.0100 | 9.7 | 20.0 |
| N-Propylbenzene | Ave | 1.066 | 1.115 | | 0.0105 | 0.0100 | 4.6 | 20.0 |
| 2-Chlorotoluene | Ave | 0.9469 | 0.9711 | | 0.0103 | 0.0100 | 2.6 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.327 | 3.472 | | 0.0104 | 0.0100 | 4.4 | 20.0 |
| 4-Chlorotoluene | Ave | 3.395 | 3.468 | | 0.0102 | 0.0100 | 2.1 | 20.0 |
| tert-Butylbenzene | Ave | 2.694 | 2.788 | | 0.0103 | 0.0100 | 3.5 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.472 | 3.597 | | 0.0104 | 0.0100 | 3.6 | 20.0 |
| sec-Butylbenzene | Ave | 3.793 | 3.984 | | 0.0105 | 0.0100 | 5.0 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.892 | 1.908 | 0.6000 | 0.0101 | 0.0100 | 0.9 | 20.0 |
| p-Isopropyltoluene | Ave | 3.286 | 3.385 | | 0.0103 | 0.0100 | 3.0 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.957 | 1.984 | 0.5000 | 0.0101 | 0.0100 | 1.4 | 20.0 |
| n-Butylbenzene | Ave | 2.732 | 2.819 | | 0.0103 | 0.0100 | 3.2 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.805 | 1.831 | 0.4000 | 0.0101 | 0.0100 | 1.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-210738/14 Calibration Date: 12/14/2015 15:24

Instrument ID: A3UX15 Calib Start Date: 12/14/2015 13:08

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 12/14/2015 15:01

Lab File ID: UXC5228.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0985 | 0.1018 | 0.0500 | 0.0103 | 0.0100 | 3.4 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 1.146 | 1.064 | 0.2000 | 0.00929 | 0.0100 | -7.1 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5163 | 0.4478 | | 0.00867 | 0.0100 | -13.3 | 20.0 |
| Naphthalene | Ave | 2.355 | 2.243 | | 0.00953 | 0.0100 | -4.7 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 1.021 | 0.8954 | | 0.00877 | 0.0100 | -12.3 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2853 | 0.2802 | | 0.00720 | 0.00733 | -1.8 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.3505 | 0.3351 | | 0.00701 | 0.00733 | -4.4 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.659 | 1.668 | | 0.00737 | 0.00733 | 0.5 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 0.6121 | 0.6703 | | 0.00803 | 0.00733 | 9.5 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-210738/15 Calibration Date: 12/14/2015 18:03

Instrument ID: A3UX15 Calib Start Date: 12/14/2015 15:47

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 12/14/2015 17:40

Lab File ID: UXC5235.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0321 | 0.0291 | | 0.0908 | 0.100 | -9.2 | 20.0 |
| Diisopropyl ether | Ave | 0.3031 | 0.3076 | | 0.0101 | 0.0100 | 1.5 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.5558 | 0.5348 | | 0.00962 | 0.0100 | -3.8 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.9024 | 0.8852 | | 0.00981 | 0.0100 | -1.9 | 20.0 |
| Ethyl acetate | Ave | 0.2030 | 0.1962 | | 0.0193 | 0.0200 | -3.3 | 20.0 |
| Propionitrile | Ave | 0.0342 | 0.0338 | | 0.0987 | 0.100 | -1.3 | 20.0 |
| Methacrylonitrile | Ave | 0.1576 | 0.1544 | | 0.0980 | 0.100 | -2.0 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.7253 | 0.7045 | | 0.00971 | 0.0100 | -2.9 | 20.0 |
| n-Butanol | Ave | 0.0046 | 0.0040 | | 0.221 | 0.250 | -11.6 | 20.0 |
| Methyl methacrylate | Ave | 0.1997 | 0.2008 | | 0.0201 | 0.0200 | 0.6 | 20.0 |
| 2-Nitropropane | Qua | | 0.0262 | | 0.0169 | 0.0200 | -15.3 | 20.0 |
| 1-Chlorohexane | Ave | 0.5469 | 0.5335 | | 0.00976 | 0.0100 | -2.4 | 20.0 |
| Cyclohexanone | Ave | 0.0165 | 0.0137 | | 0.0826 | 0.100 | -17.4 | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.571 | 3.515 | | 0.00984 | 0.0100 | -1.6 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.249 | 1.151 | | 0.00921 | 0.0100 | -7.9 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.475 | 0.7358 | | 0.00998 | 0.0200 | -50.1* | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCVIS 240-212893/2 Calibration Date: 01/04/2016 09:55

Instrument ID: A3UX15 Calib Start Date: 12/14/2015 13:08

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 12/14/2015 15:01

Lab File ID: UXC5592.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.4073 | 0.4230 | 0.1000 | 0.0104 | 0.0100 | 3.8 | 20.0 |
| Chloromethane | Ave | 0.4706 | 0.5100 | 0.1000 | 0.0108 | 0.0100 | 8.4 | 20.0 |
| Vinyl chloride | Ave | 0.4119 | 0.4332 | 0.1000 | 0.0105 | 0.0100 | 5.2 | 20.0 |
| Butadiene | Ave | 0.3871 | 0.4088 | | 0.0106 | 0.0100 | 5.6 | 20.0 |
| Bromomethane | Ave | 0.1391 | 0.1208 | 0.0500 | 0.00868 | 0.0100 | -13.2 | 20.0 |
| Chloroethane | Ave | 0.2224 | 0.2077 | 0.0500 | 0.00934 | 0.0100 | -6.6 | 20.0 |
| Dichlorofluoromethane | Ave | 0.4837 | 0.4906 | | 0.0101 | 0.0100 | 1.4 | 20.0 |
| Trichlorofluoromethane | Ave | 0.3431 | 0.3979 | 0.1000 | 0.0116 | 0.0100 | 16.0 | 20.0 |
| Ethyl ether | Ave | 0.2665 | 0.2419 | | 0.00908 | 0.0100 | -9.2 | 20.0 |
| Acrolein | Ave | 0.0381 | 0.0379 | | 0.0498 | 0.0500 | -0.5 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.3270 | 0.2919 | 0.1000 | 0.00893 | 0.0100 | -10.7 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2436 | 0.2391 | 0.0500 | 0.00982 | 0.0100 | -1.8 | 20.0 |
| Acetone | Lin1 | | 0.0781 | 0.0500 | 0.0207 | 0.0200 | 3.4 | 20.0 |
| Iodomethane | Ave | 0.5420 | 0.4830 | | 0.00891 | 0.0100 | -10.9 | 20.0 |
| Carbon disulfide | Ave | 1.048 | 0.8984 | 0.1000 | 0.00857 | 0.0100 | -14.3 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.2047 | 0.1730 | | 0.00845 | 0.0100 | -15.5 | 20.0 |
| Methyl acetate | Ave | 0.1793 | 0.1636 | 0.1000 | 0.0456 | 0.0500 | -8.7 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3487 | 0.1000 | 0.00811 | 0.0100 | -18.9 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0157 | 0.0100 | | 0.0637 | 0.100 | -36.3* | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7917 | 0.7543 | 0.1000 | 0.00953 | 0.0100 | -4.7 | 20.0 |
| Acrylonitrile | Ave | 0.0963 | 0.0937 | | 0.0973 | 0.100 | -2.7 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.3554 | 0.3176 | 0.1000 | 0.00894 | 0.0100 | -10.6 | 20.0 |
| Hexane | Ave | 0.0854 | 0.0765 | | 0.00895 | 0.0100 | -10.5 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.6528 | 0.5907 | 0.1000 | 0.00905 | 0.0100 | -9.5 | 20.0 |
| Vinyl acetate | Ave | 0.4853 | 0.3426 | | 0.00706 | 0.0100 | -29.4* | 20.0 |
| 2,2-Dichloropropane | Ave | 0.3625 | 0.2742 | | 0.00756 | 0.0100 | -24.4* | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3789 | 0.3486 | 0.1000 | 0.00920 | 0.0100 | -8.0 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.1014 | 0.1163 | 0.0500 | 0.0230 | 0.0200 | 14.8 | 20.0 |
| Bromochloromethane | Ave | 0.1659 | 0.1578 | | 0.00951 | 0.0100 | -4.9 | 20.0 |
| Tetrahydrofuran | Lin1 | | 0.0805 | | 0.0243 | 0.0200 | 21.3* | 20.0 |
| Chloroform | Ave | 0.5952 | 0.5373 | 0.2000 | 0.00903 | 0.0100 | -9.7 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.4301 | 0.3432 | 0.1000 | 0.00798 | 0.0100 | -20.2* | 20.0 |
| Cyclohexane | Ave | 0.6194 | 0.5697 | 0.1000 | 0.00920 | 0.0100 | -8.0 | 20.0 |
| Carbon tetrachloride | Ave | 0.3559 | 0.2394 | 0.1000 | 0.00673 | 0.0100 | -32.7* | 20.0 |
| 1,1-Dichloropropene | Ave | 0.4709 | 0.4290 | | 0.00911 | 0.0100 | -8.9 | 20.0 |
| Isobutyl alcohol | Ave | 0.0062 | 0.0035 | | 0.141 | 0.250 | -43.7* | 20.0 |
| Benzene | Ave | 1.528 | 1.358 | 0.5000 | 0.00888 | 0.0100 | -11.2 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.4436 | 0.4061 | 0.1000 | 0.00915 | 0.0100 | -8.5 | 20.0 |
| n-Heptane | Ave | 0.0794 | 0.0717 | | 0.00903 | 0.0100 | -9.7 | 20.0 |
| Trichloroethene | Ave | 0.3600 | 0.3283 | 0.2000 | 0.00912 | 0.0100 | -8.8 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCVIS 240-212893/2  Calibration Date: 01/04/2016 09:55

Instrument ID: A3UX15  Calib Start Date: 12/14/2015 13:08

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 12/14/2015 15:01

Lab File ID: UXC5592.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.5193 | 0.4807 | 0.1000 | 0.00926 | 0.0100 | -7.4 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.3511 | 0.3295 | 0.1000 | 0.00939 | 0.0100 | -6.1 | 20.0 |
| Dibromomethane | Ave | 0.1700 | 0.1624 | | 0.00955 | 0.0100 | -4.5 | 20.0 |
| 1,4-Dioxane | Ave | 0.0025 | 0.0026 | | 0.203 | 0.200 | 1.6 | 20.0 |
| Bromodichloromethane | Ave | 0.3598 | 0.2985 | 0.1000 | 0.00830 | 0.0100 | -17.0 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.1716 | 0.1788 | | 0.0208 | 0.0200 | 4.2 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.4575 | 0.4043 | 0.1500 | 0.00884 | 0.0100 | -11.6 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.1994 | 0.2061 | 0.0500 | 0.0207 | 0.0200 | 3.3 | 20.0 |
| Toluene | Ave | 2.026 | 1.802 | 0.4000 | 0.00890 | 0.0100 | -11.0 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.4684 | 0.3917 | 0.1000 | 0.00836 | 0.0100 | -16.4 | 20.0 |
| Ethyl methacrylate | Ave | 0.4144 | 0.3573 | | 0.00862 | 0.0100 | -13.8 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.3260 | 0.3183 | 0.1000 | 0.00976 | 0.0100 | -2.4 | 20.0 |
| Tetrachloroethene | Ave | 0.3756 | 0.3214 | 0.2000 | 0.00856 | 0.0100 | -14.4 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.6566 | 0.6173 | | 0.00940 | 0.0100 | -6.0 | 20.0 |
| 2-Hexanone | Ave | 0.1839 | 0.1857 | 0.0500 | 0.0202 | 0.0200 | 1.0 | 20.0 |
| Dibromochloromethane | Ave | 0.2904 | 0.2083 | | 0.00717 | 0.0100 | -28.3* | 20.0 |
| 1,2-Dibromoethane | Ave | 0.3299 | 0.3105 | | 0.00941 | 0.0100 | -5.9 | 20.0 |
| Chlorobenzene | Ave | 1.332 | 1.205 | 0.3000 | 0.00905 | 0.0100 | -9.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3658 | 0.2700 | | 0.00738 | 0.0100 | -26.2* | 20.0 |
| Ethylbenzene | Ave | 0.7037 | 0.6445 | | 0.00916 | 0.0100 | -8.4 | 20.0 |
| m-Xylene & p-Xylene | Ave | 1.727 | 1.549 | | 0.00897 | 0.0100 | -10.3 | 20.0 |
| o-Xylene | Ave | 0.8464 | 0.7714 | | 0.00911 | 0.0100 | -8.9 | 20.0 |
| Styrene | Ave | 1.443 | 1.355 | 0.3000 | 0.00939 | 0.0100 | -6.1 | 20.0 |
| Bromoform | Lin1 | | 0.1008 | 0.1000 | 0.00589 | 0.0100 | -41.1* | 20.0 |
| Isopropylbenzene | Ave | 2.106 | 1.910 | 0.1000 | 0.00907 | 0.0100 | -9.3 | 20.0 |
| Bromobenzene | Ave | 0.9813 | 0.8899 | | 0.00907 | 0.0100 | -9.3 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.7457 | 0.7234 | 0.3000 | 0.00970 | 0.0100 | -3.0 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.2324 | 0.2358 | | 0.0101 | 0.0100 | 1.5 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2172 | 0.1952 | | 0.00898 | 0.0100 | -10.2 | 20.0 |
| N-Propylbenzene | Ave | 1.066 | 0.9833 | | 0.00923 | 0.0100 | -7.7 | 20.0 |
| 2-Chlorotoluene | Ave | 0.9469 | 0.8482 | | 0.00896 | 0.0100 | -10.4 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.327 | 2.976 | | 0.00895 | 0.0100 | -10.5 | 20.0 |
| 4-Chlorotoluene | Ave | 3.395 | 3.006 | | 0.00885 | 0.0100 | -11.5 | 20.0 |
| tert-Butylbenzene | Ave | 2.694 | 2.448 | | 0.00909 | 0.0100 | -9.1 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.472 | 3.127 | | 0.00901 | 0.0100 | -9.9 | 20.0 |
| sec-Butylbenzene | Ave | 3.793 | 3.336 | | 0.00879 | 0.0100 | -12.1 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.892 | 1.675 | 0.6000 | 0.00885 | 0.0100 | -11.5 | 20.0 |
| p-Isopropyltoluene | Ave | 3.286 | 2.914 | | 0.00887 | 0.0100 | -11.3 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.957 | 1.714 | 0.5000 | 0.00876 | 0.0100 | -12.4 | 20.0 |
| n-Butylbenzene | Ave | 2.732 | 2.311 | | 0.00846 | 0.0100 | -15.4 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.805 | 1.604 | 0.4000 | 0.00889 | 0.0100 | -11.1 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCVIS 240-212893/2  Calibration Date: 01/04/2016 09:55

Instrument ID: A3UX15  Calib Start Date: 12/14/2015 13:08

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 12/14/2015 15:01

Lab File ID: UXC5592.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0985 | 0.0717 | 0.0500 | 0.00728 | 0.0100 | -27.2* | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 1.146 | 0.9020 | 0.2000 | 0.00787 | 0.0100 | -21.3* | 20.0 |
| Hexachlorobutadiene | Ave | 0.5163 | 0.3423 | | 0.00663 | 0.0100 | -33.7* | 20.0 |
| Naphthalene | Ave | 2.355 | 2.014 | | 0.00855 | 0.0100 | -14.5 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 1.021 | 0.7653 | | 0.00749 | 0.0100 | -25.1* | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2853 | 0.2864 | | 0.00736 | 0.00733 | 0.4 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.3505 | 0.3364 | | 0.00704 | 0.00733 | -4.0 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.659 | 1.672 | | 0.00739 | 0.00733 | 0.8 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 0.6121 | 0.6848 | | 0.00820 | 0.00733 | 11.9 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212893/3 Calibration Date: 01/04/2016 10:41

Instrument ID: A3UX15 Calib Start Date: 12/14/2015 15:47

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 12/14/2015 17:40

Lab File ID: UXC5594.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0321 | 0.0282 | | 0.0878 | 0.100 | -12.2 | 20.0 |
| Diisopropyl ether | Ave | 0.3031 | 0.2927 | | 0.00966 | 0.0100 | -3.4 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.5558 | 0.4890 | | 0.00880 | 0.0100 | -12.0 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.9024 | 0.8266 | | 0.00916 | 0.0100 | -8.4 | 20.0 |
| Ethyl acetate | Ave | 0.2030 | 0.1864 | | 0.0184 | 0.0200 | -8.2 | 20.0 |
| Propionitrile | Ave | 0.0342 | 0.0329 | | 0.0963 | 0.100 | -3.7 | 20.0 |
| Methacrylonitrile | Ave | 0.1576 | 0.1494 | | 0.0948 | 0.100 | -5.2 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.7253 | 0.6808 | | 0.00939 | 0.0100 | -6.1 | 20.0 |
| n-Butanol | Ave | 0.0046 | 0.0029 | | 0.159 | 0.250 | -36.5* | 20.0 |
| Methyl methacrylate | Ave | 0.1997 | 0.1890 | | 0.0189 | 0.0200 | -5.3 | 20.0 |
| 2-Nitropropane | Qua | | 0.0143 | | 0.0102 | 0.0200 | -49.2* | 20.0 |
| 1-Chlorohexane | Ave | 0.5469 | 0.4819 | | 0.00881 | 0.0100 | -11.9 | 20.0 |
| Cyclohexanone | Ave | 0.0165 | 0.0146 | | 0.0885 | 0.100 | -11.5 | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.571 | 3.251 | | 0.00910 | 0.0100 | -9.0 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.249 | 1.057 | | 0.00846 | 0.0100 | -15.4 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.475 | 0.6915 | | 0.00938 | 0.0200 | -53.1* | 20.0 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212606/24

Matrix: Solid | Lab File ID: 147321.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 12/30/2015 20:00

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.80 | U | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.562 | J | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.78 | U | 5.0 | 0.78 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.460 | J | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 1.1 | U | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.90 | U | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212606/24

Matrix: Solid | Lab File ID: 147321.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 12/30/2015 20:00

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.79 | U | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.71 | U | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.7 | U | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.74 | U | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.618 | J | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.91 | U | 5.0 | 0.91 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.80 | U | 5.0 | 0.80 |
| 108-88-3 | Toluene | 0.27 | U | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212606/24

Matrix: Solid | Lab File ID: 147321.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 12/30/2015 20:00

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 88 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 93 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 86 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 97 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212893/8

Matrix: Water | Lab File ID: UXC5596.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 11:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.28 | U | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.44 | U | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.22 | U | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.30 | U | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 0.45 | U | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.37 | U | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.44 | U | 1.0 | 0.44 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.47 | U | 5.0 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.32 | U | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.41 | U | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.82 | U | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.25 | U | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 0.23 | U | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 0.25 | U | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.19 | U | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 0.19 | U | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.27 | U | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 0.26 | U | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 1.34 | J | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.59 | U | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 0.40 | U | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 0.48 | U | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 10 | 0.99 |
| 67-64-1 | Acetone | 0.94 | U | 10 | 0.94 |
| 71-43-2 | Benzene | 0.35 | U | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 0.35 | U | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 0.56 | U | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 0.44 | U | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.38 | U | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 0.43 | U | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 0.25 | U | 1.0 | 0.25 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212893/8

Matrix: Water | Lab File ID: UXC5596.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 11:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 1.0 | 0.43 |
| 75-00-3 | Chloroethane | 0.32 | U | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 0.25 | U | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 0.44 | U | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.26 | U | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.46 | U | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 0.42 | U | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.50 | U | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 4.9 | U | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 0.25 | U | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 0.35 | U | 1.0 | 0.35 |
| 108-20-3 | Diisopropyl ether | 0.50 | U | 10 | 0.50 |
| 98-82-8 | Isopropylbenzene | 0.35 | U | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 0.20 | U | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 0.33 | U | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.24 | U | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 0.45 | U | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 0.31 | U | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 0.40 | U | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 0.25 | U | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 0.48 | U | 1.0 | 0.48 |
| 100-42-5 | Styrene | 0.45 | U | 1.0 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.30 | U | 5.0 | 0.30 |
| 637-92-3 | Ethyl t-butyl ether | 0.23 | U | 5.0 | 0.23 |
| 127-18-4 | Tetrachloroethene | 0.31 | U | 1.0 | 0.31 |
| 108-88-3 | Toluene | 0.23 | U | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.30 | U | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.56 | U | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 0.22 | U | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 0.49 | U | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 0.41 | U | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 0.29 | U | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 0.52 | U | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 0.41 | U | 1.0 | 0.41 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212893/8

Matrix: Water | Lab File ID: UXC5596.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 11:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.32 | U | 1.0 | 0.32 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 102 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 94 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 111 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 97 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212606/23

Matrix: Solid | Lab File ID: 147320.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 12/30/2015 19:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 58.6 | | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 54.0 | | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 53.5 | | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 45.3 | | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 51.1 | | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 42.2 | | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 54.6 | | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 45.6 | | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 57.2 | | 5.0 | 0.78 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 45.6 | | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 52.6 | | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 55.6 | | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 48.1 | | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 52.5 | | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 52.7 | | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 47.0 | | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 51.7 | | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 47.3 | | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 54.3 | | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 96.3 | | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 54.3 | | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 52.2 | | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 99.2 | | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 48.3 | | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 53.4 | | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 109 | | 20 | 1.2 |
| 67-64-1 | Acetone | 94.9 | | 20 | 5.7 |
| 71-43-2 | Benzene | 50.8 | | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 51.6 | | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 54.9 | | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 19.2 | | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 54.9 | | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 57.7 | | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 50.5 | | 5.0 | 0.53 |
| 124-48-1 | Dibromochloromethane | 54.7 | | 5.0 | 0.37 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212606/23

Matrix: Solid | Lab File ID: 147320.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 12/30/2015 19:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 19.5 | | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 51.8 | | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 20.5 | | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 52.6 | | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 53.5 | | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 53.8 | | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 52.0 | | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 20.7 | | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 54.4 | | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 450 | | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 51.6 | | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 51.3 | | 5.0 | 0.40 |
| 98-82-8 | Isopropylbenzene | 54.4 | | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 53.0 | | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 50.7 | | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 53.8 | | 10 | 0.42 |
| 91-20-3 | Naphthalene | 51.1 | | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 50.3 | | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 53.2 | | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 54.6 | | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 55.3 | | 5.0 | 0.91 |
| 100-42-5 | Styrene | 53.8 | | 5.0 | 0.40 |
| 127-18-4 | Tetrachloroethene | 51.4 | | 5.0 | 0.80 |
| 108-88-3 | Toluene | 53.2 | | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 53.2 | | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 55.5 | | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 53.1 | | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 21.1 | | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 48.6 | | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 20.2 | | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 108 | | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 54.9 | | 5.0 | 0.67 |
| 106-93-4 | 1,2-Dibromoethane | 56.1 | | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-212606/23

Matrix: Solid  Lab File ID: 147320.D

Analysis Method: 8260C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 12/30/2015 19:39

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: DB-624 ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 212606  Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 97 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 94 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 91 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 92 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212606/25

Matrix: Solid | Lab File ID: 147322.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 12/30/2015 20:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 526-73-8 | 1,2,3-Trimethylbenzene | 53.4 | | 5.0 | 0.46 |
| 108-20-3 | Diisopropyl ether | 55.3 | | 10 | 0.59 |
| 994-05-8 | Tert-amyl methyl ether | 51.1 | | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 52.6 | | 5.0 | 0.41 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 93 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 90 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 89 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 92 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212893/5

Matrix: Water | Lab File ID: UXC5593.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 10:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 8.23 | | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 8.79 | | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 9.40 | | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 10.5 | | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 9.36 | | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 9.64 | | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 10.1 | | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 8.36 | | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 9.73 | | 1.0 | 0.44 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 8.53 | | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 9.27 | | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 8.17 | | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 9.02 | | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 9.32 | | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 9.80 | | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 9.06 | | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 9.36 | | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 9.07 | | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 7.92 | | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 21.5 | | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 9.80 | J | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 9.14 | | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 18.5 | | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 9.06 | | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 9.23 | | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 19.6 | | 10 | 0.99 |
| 67-64-1 | Acetone | 19.1 | | 10 | 0.94 |
| 71-43-2 | Benzene | 9.39 | | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 9.14 | | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 6.29 | | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 6.43 | | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 9.43 | | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 7.89 | | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 9.32 | | 1.0 | 0.25 |
| 124-48-1 | Dibromochloromethane | 7.97 | | 1.0 | 0.43 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212893/5

Matrix: Water | Lab File ID: UXC5593.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 10:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 75-00-3 | Chloroethane | 7.54 | | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 9.57 | | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 9.55 | | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 9.75 | | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 9.84 | | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 9.81 | | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 9.83 | | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 7.57 | | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 8.91 | | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 79.5 | | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 9.45 | | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 7.63 | | 1.0 | 0.35 |
| 98-82-8 | Isopropylbenzene | 9.47 | | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 9.50 | | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 8.56 | | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 9.38 | | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 9.11 | | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 8.99 | | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 9.29 | | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 9.59 | | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 9.27 | | 1.0 | 0.48 |
| 100-42-5 | Styrene | 9.62 | | 1.0 | 0.45 |
| 127-18-4 | Tetrachloroethene | 9.39 | | 1.0 | 0.31 |
| 108-88-3 | Toluene | 9.29 | | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 9.74 | | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 9.12 | | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 9.82 | | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 9.18 | | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 8.16 | | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 9.13 | | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 19.0 | | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 9.36 | | 1.0 | 0.41 |
| 106-93-4 | 1,2-Dibromoethane | 9.65 | | 1.0 | 0.32 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Client Sample ID: Lab Sample ID: LCS 240-212893/5

Matrix: Water Lab File ID: UXC5593.D

Analysis Method: 8260C Date Collected:

Sample wt/vol: 5(mL) Date Analyzed: 01/04/2016 10:18

Soil Aliquot Vol: Dilution Factor: 1

Soil Extract Vol.: GC Column: DB-624 ID: 0.18(mm)

% Moisture: Level: (low/med) Low

Analysis Batch No.: 212893 Units: ug/L

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 100 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 114 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 96 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MS | Lab Sample ID: 240-59482-2 MS

Matrix: Solid | Lab File ID: 147327.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.65(g) | Date Analyzed: 12/30/2015 22:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 44.5 | | 5.4 | 0.54 |
| 71-55-6 | 1,1,1-Trichloroethane | 50.6 | | 5.4 | 0.71 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 43.6 | | 5.4 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 52.2 | | 5.4 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 45.9 | | 5.4 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 48.2 | | 5.4 | 0.87 |
| 563-58-6 | 1,1-Dichloropropene | 50.4 | | 5.4 | 0.40 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 22.8 | | 5.4 | 0.54 |
| 96-18-4 | 1,2,3-Trichloropropane | 48.1 | | 5.4 | 0.85 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.789 | J | 5.4 | 0.50 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 23.3 | | 5.4 | 0.41 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 37.0 | | 5.4 | 0.63 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 40.1 | | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 30.9 | | 5.4 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 44.3 | | 5.4 | 0.50 |
| 78-87-5 | 1,2-Dichloropropane | 47.0 | | 5.4 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 29.9 | | 5.4 | 0.75 |
| 142-28-9 | 1,3-Dichloropropane | 41.4 | | 5.4 | 0.43 |
| 106-46-7 | 1,4-Dichlorobenzene | 27.7 | | 5.4 | 0.30 |
| 594-20-7 | 2,2-Dichloropropane | 46.0 | | 5.4 | 0.48 |
| 78-93-3 | 2-Butanone (MEK) | 89.2 | | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 33.0 | J | 54 | 0.69 |
| 95-49-8 | 2-Chlorotoluene | 35.4 | | 5.4 | 0.33 |
| 591-78-6 | 2-Hexanone | 90.1 | | 22 | 0.74 |
| 106-43-4 | 4-Chlorotoluene | 32.1 | | 5.4 | 0.28 |
| 99-87-6 | p-Isopropyltoluene | 34.2 | | 5.4 | 0.44 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 97.1 | | 22 | 1.3 |
| 67-64-1 | Acetone | 102 | | 22 | 6.2 |
| 71-43-2 | Benzene | 45.9 | | 5.4 | 0.98 |
| 108-86-1 | Bromobenzene | 35.4 | | 5.4 | 0.36 |
| 75-25-2 | Bromoform | 42.3 | | 5.4 | 0.26 |
| 74-83-9 | Bromomethane | 16.0 | | 5.4 | 0.46 |
| 75-15-0 | Carbon disulfide | 43.4 | | 5.4 | 0.64 |
| 56-23-5 | Carbon tetrachloride | 54.2 | | 5.4 | 0.71 |
| 108-90-7 | Chlorobenzene | 38.3 | | 5.4 | 0.58 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MS | Lab Sample ID: 240-59482-2 MS

Matrix: Solid | Lab File ID: 147327.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.65(g) | Date Analyzed: 12/30/2015 22:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 42.4 | | 5.4 | 0.40 |
| 75-00-3 | Chloroethane | 17.2 | | 5.4 | 0.48 |
| 67-66-3 | Chloroform | 45.3 | | 5.4 | 0.39 |
| 74-87-3 | Chloromethane | 18.5 | | 5.4 | 0.86 |
| 156-59-2 | cis-1,2-Dichloroethene | 43.2 | | 5.4 | 0.30 |
| 10061-01-5 | cis-1,3-Dichloropropene | 42.5 | | 5.4 | 0.77 |
| 74-95-3 | Dibromomethane | 43.3 | | 5.4 | 0.46 |
| 74-97-5 | Bromochloromethane | 41.7 | | 5.4 | 0.36 |
| 75-71-8 | Dichlorodifluoromethane | 20.5 | | 5.4 | 0.30 |
| 75-27-4 | Bromodichloromethane | 46.2 | | 5.4 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 454 | | 220 | 6.2 |
| 100-41-4 | Ethylbenzene | 40.5 | | 5.4 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 24.2 | | 5.4 | 0.43 |
| 108-20-3 | Diisopropyl ether | 0.64 | U | 11 | 0.64 |
| 98-82-8 | Isopropylbenzene | 39.5 | | 5.4 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 46.4 | | 5.4 | 0.47 |
| 75-09-2 | Methylene Chloride | 40.6 | | 5.4 | 0.80 |
| 179601-23-1 | m-Xylene & p-Xylene | 42.1 | | 11 | 0.46 |
| 91-20-3 | Naphthalene | 31.1 | | 5.4 | 0.38 |
| 104-51-8 | n-Butylbenzene | 29.5 | | 5.4 | 0.64 |
| 103-65-1 | N-Propylbenzene | 36.8 | | 5.4 | 0.65 |
| 95-47-6 | o-Xylene | 42.2 | | 5.4 | 0.62 |
| 135-98-8 | sec-Butylbenzene | 35.7 | | 5.4 | 0.99 |
| 100-42-5 | Styrene | 37.2 | | 5.4 | 0.43 |
| 994-05-8 | Tert-amyl methyl ether | 0.47 | U | 5.4 | 0.47 |
| 637-92-3 | Ethyl t-butyl ether | 0.44 | U | 5.4 | 0.44 |
| 127-18-4 | Tetrachloroethene | 42.6 | | 5.4 | 0.87 |
| 108-88-3 | Toluene | 45.1 | | 5.4 | 0.29 |
| 156-60-5 | trans-1,2-Dichloroethene | 44.0 | | 5.4 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 37.3 | | 5.4 | 0.31 |
| 79-01-6 | Trichloroethene | 44.7 | | 5.4 | 0.41 |
| 75-69-4 | Trichlorofluoromethane | 20.8 | | 5.4 | 0.31 |
| 108-05-4 | Vinyl acetate | 32.9 | | 11 | 0.72 |
| 75-01-4 | Vinyl chloride | 18.3 | | 5.4 | 0.33 |
| 1330-20-7 | Xylenes, Total | 84.3 | | 11 | 0.59 |
| 98-06-6 | tert-Butylbenzene | 38.2 | | 5.4 | 0.73 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MS | Lab Sample ID: 240-59482-2 MS

Matrix: Solid | Lab File ID: 147327.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.65(g) | Date Analyzed: 12/30/2015 22:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 42.6 | | 5.4 | 0.58 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 92 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 95 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 91 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 97 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-59545-B-2 MS

Matrix: Water | Lab File ID: UXC5610.D

Analysis Method: 8260C | Date Collected: 12/29/2015 16:48

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 17:00

Soil Aliquot Vol: | Dilution Factor: 500

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 71-55-6 | 1,1,1-Trichloroethane | 4250 | | 500 | 220 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 4970 | | 500 | 110 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 4400 | | 500 | 230 |
| 75-34-3 | 1,1-Dichloroethane | 10600 | | 500 | 150 |
| 75-35-4 | 1,1-Dichloroethene | 4740 | | 500 | 230 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 3990 | | 500 | 160 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3890 | | 1000 | 410 |
| 95-50-1 | 1,2-Dichlorobenzene | 4420 | | 500 | 130 |
| 107-06-2 | 1,2-Dichloroethane | 4970 | | 500 | 120 |
| 78-87-5 | 1,2-Dichloropropane | 4850 | | 500 | 130 |
| 541-73-1 | 1,3-Dichlorobenzene | 4440 | | 500 | 95 |
| 106-46-7 | 1,4-Dichlorobenzene | 4470 | | 500 | 140 |
| 78-93-3 | 2-Butanone (MEK) | 11300 | | 5000 | 270 |
| 591-78-6 | 2-Hexanone | 10100 | | 5000 | 240 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 10400 | | 5000 | 500 |
| 67-64-1 | Acetone | 8410 | | 5000 | 470 |
| 71-43-2 | Benzene | 4720 | | 500 | 180 |
| 75-25-2 | Bromoform | 3020 | | 500 | 280 |
| 74-83-9 | Bromomethane | 2680 | | 500 | 220 |
| 75-15-0 | Carbon disulfide | 4490 | | 500 | 190 |
| 56-23-5 | Carbon tetrachloride | 3710 | | 500 | 220 |
| 108-90-7 | Chlorobenzene | 4670 | | 500 | 130 |
| 124-48-1 | Dibromochloromethane | 3910 | | 500 | 220 |
| 75-00-3 | Chloroethane | 3500 | | 500 | 160 |
| 67-66-3 | Chloroform | 4760 | | 500 | 130 |
| 74-87-3 | Chloromethane | 4750 | | 500 | 220 |
| 156-59-2 | cis-1,2-Dichloroethene | 17700 | | 500 | 130 |
| 10061-01-5 | cis-1,3-Dichloropropene | 4550 | | 500 | 230 |
| 75-71-8 | Dichlorodifluoromethane | 3190 | | 500 | 160 |
| 75-27-4 | Bromodichloromethane | 4260 | | 500 | 150 |
| 100-41-4 | Ethylbenzene | 4710 | | 500 | 130 |
| 98-82-8 | Isopropylbenzene | 4710 | | 500 | 180 |
| 1634-04-4 | Methyl tert-butyl ether | 4640 | | 500 | 100 |
| 75-09-2 | Methylene Chloride | 4360 | | 500 | 170 |
| 179601-23-1 | m-Xylene & p-Xylene | 4700 | | 1000 | 120 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-59545-B-2 MS

Matrix: Water | Lab File ID: UXC5610.D

Analysis Method: 8260C | Date Collected: 12/29/2015 16:48

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 17:00

Soil Aliquot Vol: | Dilution Factor: 500

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 95-47-6 | o-Xylene | 4760 | | 500 | 130 |
| 100-42-5 | Styrene | 4790 | | 500 | 230 |
| 127-18-4 | Tetrachloroethene | 4540 | | 500 | 160 |
| 108-88-3 | Toluene | 4690 | | 500 | 120 |
| 156-60-5 | trans-1,2-Dichloroethene | 5390 | | 500 | 150 |
| 10061-02-6 | trans-1,3-Dichloropropene | 4330 | | 500 | 280 |
| 79-01-6 | Trichloroethene | 4800 | | 500 | 110 |
| 75-69-4 | Trichlorofluoromethane | 3960 | | 500 | 250 |
| 75-01-4 | Vinyl chloride | 8920 | | 500 | 150 |
| 1330-20-7 | Xylenes, Total | 9460 | | 1000 | 260 |
| 106-93-4 | 1,2-Dibromoethane | 5040 | | 500 | 160 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 103 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 100 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 117 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MSD | Lab Sample ID: 240-59482-2 MSD

Matrix: Solid | Lab File ID: 147328.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.47(g) | Date Analyzed: 12/30/2015 22:31

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 41.7 | | 5.6 | 0.56 |
| 71-55-6 | 1,1,1-Trichloroethane | 55.0 | | 5.6 | 0.73 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 47.6 | | 5.6 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 58.3 | | 5.6 | 0.47 |
| 75-34-3 | 1,1-Dichloroethane | 50.9 | | 5.6 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 56.9 | | 5.6 | 0.90 |
| 563-58-6 | 1,1-Dichloropropene | 51.9 | | 5.6 | 0.41 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 19.0 | | 5.6 | 0.56 |
| 96-18-4 | 1,2,3-Trichloropropane | 50.5 | | 5.6 | 0.87 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 1.61 | J | 5.6 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 17.4 | | 5.6 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 27.3 | | 5.6 | 0.65 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 48.9 | | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 23.5 | | 5.6 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 49.6 | | 5.6 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 48.5 | | 5.6 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 20.8 | | 5.6 | 0.77 |
| 142-28-9 | 1,3-Dichloropropane | 43.5 | | 5.6 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 20.0 | | 5.6 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 50.4 | | 5.6 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 106 | | 22 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 39.9 | J | 56 | 0.72 |
| 95-49-8 | 2-Chlorotoluene | 24.1 | | 5.6 | 0.34 |
| 591-78-6 | 2-Hexanone | 103 | | 22 | 0.76 |
| 106-43-4 | 4-Chlorotoluene | 21.3 | | 5.6 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 20.0 | | 5.6 | 0.46 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 117 | | 22 | 1.4 |
| 67-64-1 | Acetone | 125 | | 22 | 6.4 |
| 71-43-2 | Benzene | 47.6 | | 5.6 | 1.0 |
| 108-86-1 | Bromobenzene | 28.4 | | 5.6 | 0.37 |
| 75-25-2 | Bromoform | 42.8 | | 5.6 | 0.27 |
| 74-83-9 | Bromomethane | 17.9 | | 5.6 | 0.47 |
| 75-15-0 | Carbon disulfide | 45.7 | | 5.6 | 0.66 |
| 56-23-5 | Carbon tetrachloride | 56.7 | | 5.6 | 0.73 |
| 108-90-7 | Chlorobenzene | 31.6 | | 5.6 | 0.59 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MSD | Lab Sample ID: 240-59482-2 MSD

Matrix: Solid | Lab File ID: 147328.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.47(g) | Date Analyzed: 12/30/2015 22:31

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 43.6 | | 5.6 | 0.41 |
| 75-00-3 | Chloroethane | 19.1 | | 5.6 | 0.49 |
| 67-66-3 | Chloroform | 49.0 | | 5.6 | 0.40 |
| 74-87-3 | Chloromethane | 22.6 | | 5.6 | 0.89 |
| 156-59-2 | cis-1,2-Dichloroethene | 47.7 | | 5.6 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 41.8 | | 5.6 | 0.80 |
| 74-95-3 | Dibromomethane | 46.2 | | 5.6 | 0.47 |
| 74-97-5 | Bromochloromethane | 46.8 | | 5.6 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 24.1 | | 5.6 | 0.31 |
| 75-27-4 | Bromodichloromethane | 47.1 | | 5.6 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 541 | | 220 | 6.4 |
| 100-41-4 | Ethylbenzene | 31.4 | | 5.6 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 12.2 | | 5.6 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.66 | U | 11 | 0.66 |
| 98-82-8 | Isopropylbenzene | 28.6 | | 5.6 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 53.0 | | 5.6 | 0.48 |
| 75-09-2 | Methylene Chloride | 45.3 | | 5.6 | 0.83 |
| 179601-23-1 | m-Xylene & p-Xylene | 34.9 | | 11 | 0.47 |
| 91-20-3 | Naphthalene | 31.3 | | 5.6 | 0.39 |
| 104-51-8 | n-Butylbenzene | 15.6 | | 5.6 | 0.66 |
| 103-65-1 | N-Propylbenzene | 23.4 | | 5.6 | 0.67 |
| 95-47-6 | o-Xylene | 33.9 | | 5.6 | 0.64 |
| 135-98-8 | sec-Butylbenzene | 21.4 | | 5.6 | 1.0 |
| 100-42-5 | Styrene | 28.1 | | 5.6 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.6 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.46 | U | 5.6 | 0.46 |
| 127-18-4 | Tetrachloroethene | 34.7 | | 5.6 | 0.90 |
| 108-88-3 | Toluene | 40.2 | | 5.6 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 48.3 | | 5.6 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 38.8 | | 5.6 | 0.32 |
| 79-01-6 | Trichloroethene | 44.3 | | 5.6 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 24.7 | | 5.6 | 0.32 |
| 108-05-4 | Vinyl acetate | 43.6 | | 11 | 0.74 |
| 75-01-4 | Vinyl chloride | 21.8 | | 5.6 | 0.34 |
| 1330-20-7 | Xylenes, Total | 68.8 | | 11 | 0.61 |
| 98-06-6 | tert-Butylbenzene | 25.3 | | 5.6 | 0.75 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MSD | Lab Sample ID: 240-59482-2 MSD

Matrix: Solid | Lab File ID: 147328.D

Analysis Method: 8260C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.47(g) | Date Analyzed: 12/30/2015 22:31

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212606 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 43.9 | | 5.6 | 0.59 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 91 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 95 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 100 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-59545-B-2 MSD

Matrix: Water | Lab File ID: UXC5611.D

Analysis Method: 8260C | Date Collected: 12/29/2015 16:48

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 17:23

Soil Aliquot Vol: | Dilution Factor: 500

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 71-55-6 | 1,1,1-Trichloroethane | 4480 | | 500 | 220 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 4970 | | 500 | 110 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 5180 | | 500 | 230 |
| 75-34-3 | 1,1-Dichloroethane | 10800 | | 500 | 150 |
| 75-35-4 | 1,1-Dichloroethene | 5020 | | 500 | 230 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 4280 | | 500 | 160 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3860 | | 1000 | 410 |
| 95-50-1 | 1,2-Dichlorobenzene | 4610 | | 500 | 130 |
| 107-06-2 | 1,2-Dichloroethane | 4960 | | 500 | 120 |
| 78-87-5 | 1,2-Dichloropropane | 4910 | | 500 | 130 |
| 541-73-1 | 1,3-Dichlorobenzene | 4550 | | 500 | 95 |
| 106-46-7 | 1,4-Dichlorobenzene | 4580 | | 500 | 140 |
| 78-93-3 | 2-Butanone (MEK) | 9840 | | 5000 | 270 |
| 591-78-6 | 2-Hexanone | 8950 | | 5000 | 240 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 9860 | | 5000 | 500 |
| 67-64-1 | Acetone | 8100 | | 5000 | 470 |
| 71-43-2 | Benzene | 4820 | | 500 | 180 |
| 75-25-2 | Bromoform | 3200 | | 500 | 280 |
| 74-83-9 | Bromomethane | 5100 | | 500 | 220 |
| 75-15-0 | Carbon disulfide | 4730 | | 500 | 190 |
| 56-23-5 | Carbon tetrachloride | 4200 | | 500 | 220 |
| 108-90-7 | Chlorobenzene | 4690 | | 500 | 130 |
| 124-48-1 | Dibromochloromethane | 3960 | | 500 | 220 |
| 75-00-3 | Chloroethane | 5260 | | 500 | 160 |
| 67-66-3 | Chloroform | 4950 | | 500 | 130 |
| 74-87-3 | Chloromethane | 4860 | | 500 | 220 |
| 156-59-2 | cis-1,2-Dichloroethene | 18400 | | 500 | 130 |
| 10061-01-5 | cis-1,3-Dichloropropene | 4610 | | 500 | 230 |
| 75-71-8 | Dichlorodifluoromethane | 4070 | | 500 | 160 |
| 75-27-4 | Bromodichloromethane | 4410 | | 500 | 150 |
| 100-41-4 | Ethylbenzene | 4850 | | 500 | 130 |
| 98-82-8 | Isopropylbenzene | 4950 | | 500 | 180 |
| 1634-04-4 | Methyl tert-butyl ether | 4740 | | 500 | 100 |
| 75-09-2 | Methylene Chloride | 4590 | | 500 | 170 |
| 179601-23-1 | m-Xylene & p-Xylene | 4800 | | 1000 | 120 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-59545-B-2 MSD

Matrix: Water | Lab File ID: UXC5611.D

Analysis Method: 8260C | Date Collected: 12/29/2015 16:48

Sample wt/vol: 5(mL) | Date Analyzed: 01/04/2016 17:23

Soil Aliquot Vol: | Dilution Factor: 500

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 212893 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 95-47-6 | o-Xylene | 4820 | | 500 | 130 |
| 100-42-5 | Styrene | 4900 | | 500 | 230 |
| 127-18-4 | Tetrachloroethene | 4620 | | 500 | 160 |
| 108-88-3 | Toluene | 4620 | | 500 | 120 |
| 156-60-5 | trans-1,2-Dichloroethene | 5740 | | 500 | 150 |
| 10061-02-6 | trans-1,3-Dichloropropene | 4270 | | 500 | 280 |
| 79-01-6 | Trichloroethene | 4960 | | 500 | 110 |
| 75-69-4 | Trichlorofluoromethane | 4740 | | 500 | 250 |
| 75-01-4 | Vinyl chloride | 9380 | | 500 | 150 |
| 1330-20-7 | Xylenes, Total | 9620 | | 1000 | 260 |
| 106-93-4 | 1,2-Dibromoethane | 4900 | | 500 | 160 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 98 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 105 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 116 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 97 | | 78-125 |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A3UX14 Start Date: 12/30/2015 11:47

Analysis Batch Number: 212606 End Date: 12/30/2015 23:14

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-212606/29 | | 12/30/2015 11:47 | 1 | BFB14175.D | DB-624 0.18(mm) |
| IC 240-212606/3 | | 12/30/2015 12:29 | 1 | 147300.D | DB-624 0.18(mm) |
| IC 240-212606/4 | | 12/30/2015 12:50 | 1 | 147301.D | DB-624 0.18(mm) |
| IC 240-212606/5 | | 12/30/2015 13:12 | 1 | 147302.D | DB-624 0.18(mm) |
| ICIS 240-212606/6 | | 12/30/2015 13:33 | 1 | 147303.D | DB-624 0.18(mm) |
| IC 240-212606/7 | | 12/30/2015 13:55 | 1 | 147304.D | DB-624 0.18(mm) |
| IC 240-212606/8 | | 12/30/2015 14:16 | 1 | 147305.D | DB-624 0.18(mm) |
| IC 240-212606/9 | | 12/30/2015 14:38 | 1 | 147306.D | DB-624 0.18(mm) |
| IC 240-212606/10 | | 12/30/2015 14:59 | 1 | 147307.D | DB-624 0.18(mm) |
| IC 240-212606/11 | | 12/30/2015 15:21 | 1 | 147308.D | DB-624 0.18(mm) |
| IC 240-212606/12 | | 12/30/2015 15:42 | 1 | 147309.D | DB-624 0.18(mm) |
| IC 240-212606/13 | | 12/30/2015 16:04 | 1 | 147310.D | DB-624 0.18(mm) |
| IC 240-212606/14 | | 12/30/2015 16:25 | 1 | 147311.D | DB-624 0.18(mm) |
| IC 240-212606/15 | | 12/30/2015 16:47 | 1 | 147312.D | DB-624 0.18(mm) |
| IC 240-212606/16 | | 12/30/2015 17:08 | 1 | 147313.D | DB-624 0.18(mm) |
| IC 240-212606/17 | | 12/30/2015 17:30 | 1 | 147314.D | DB-624 0.18(mm) |
| IC 240-212606/18 | | 12/30/2015 17:51 | 1 | 147315.D | DB-624 0.18(mm) |
| IC 240-212606/19 | | 12/30/2015 18:13 | 1 | 147316.D | DB-624 0.18(mm) |
| IC 240-212606/20 | | 12/30/2015 18:34 | 1 | 147317.D | DB-624 0.18(mm) |
| ICV 240-212606/21 | | 12/30/2015 18:56 | 1 | 147318.D | DB-624 0.18(mm) |
| ICV 240-212606/22 | | 12/30/2015 19:17 | 1 | 147319.D | DB-624 0.18(mm) |
| LCS 240-212606/23 | | 12/30/2015 19:39 | 1 | 147320.D | DB-624 0.18(mm) |
| MB 240-212606/24 | | 12/30/2015 20:00 | 1 | 147321.D | DB-624 0.18(mm) |
| LCS 240-212606/25 | | 12/30/2015 20:22 | 1 | 147322.D | DB-624 0.18(mm) |
| ZZZZZ | | 12/30/2015 20:43 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 12/30/2015 21:05 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 12/30/2015 21:26 | 1 | | DB-624 0.18(mm) |
| 240-59482-2 | F-VB-UST-P-7 | 12/30/2015 21:48 | 1 | 147326.D | DB-624 0.18(mm) |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 12/30/2015 22:09 | 1 | 147327.D | DB-624 0.18(mm) |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 12/30/2015 22:31 | 1 | 147328.D | DB-624 0.18(mm) |
| 240-59482-3 | F-VB-UST-P-8 | 12/30/2015 23:14 | 1 | 147330.D | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A3UX15 Start Date: 12/14/2015 12:30

Analysis Batch Number: 210738 End Date: 12/14/2015 20:40

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-210738/1 | | 12/14/2015 12:30 | 1 | BFB496.D | DB-624 0.18(mm) |
| STD8260 240-210738/2 IC | | 12/14/2015 13:08 | 1 | UXC5222A.D | DB-624 0.18(mm) |
| STD8260 240-210738/3 IC | | 12/14/2015 13:31 | 1 | UXC5223.D | DB-624 0.18(mm) |
| STD8260 240-210738/4 ICIS | | 12/14/2015 13:54 | 1 | UXC5224.D | DB-624 0.18(mm) |
| STD8260 240-210738/5 IC | | 12/14/2015 14:16 | 1 | UXC5225.D | DB-624 0.18(mm) |
| STD8260 240-210738/6 IC | | 12/14/2015 14:39 | 1 | UXC5226.D | DB-624 0.18(mm) |
| STD8260 240-210738/7 IC | | 12/14/2015 15:01 | 1 | UXC5227.D | DB-624 0.18(mm) |
| ICV 240-210738/14 | | 12/14/2015 15:24 | 1 | UXC5228.D | DB-624 0.18(mm) |
| STD 240-210738/8 IC | | 12/14/2015 15:47 | 1 | UXC5229.D | DB-624 0.18(mm) |
| STD 240-210738/9 IC | | 12/14/2015 16:09 | 1 | UXC5230.D | DB-624 0.18(mm) |
| STD 240-210738/10 IC | | 12/14/2015 16:32 | 1 | UXC5231.D | DB-624 0.18(mm) |
| STD 240-210738/11 IC | | 12/14/2015 16:55 | 1 | UXC5232.D | DB-624 0.18(mm) |
| STD 240-210738/12 IC | | 12/14/2015 17:17 | 1 | UXC5233.D | DB-624 0.18(mm) |
| STD 240-210738/13 IC | | 12/14/2015 17:40 | 1 | UXC5234.D | DB-624 0.18(mm) |
| ICV 240-210738/15 | | 12/14/2015 18:03 | 1 | UXC5235.D | DB-624 0.18(mm) |
| STDTHT 240-210738/17 IC | | 12/14/2015 18:47 | 1 | | DB-624 0.18(mm) |
| STDTHT 240-210738/18 IC | | 12/14/2015 19:10 | 1 | | DB-624 0.18(mm) |
| STDTHT 240-210738/19 IC | | 12/14/2015 19:32 | 1 | | DB-624 0.18(mm) |
| STDTHT 240-210738/20 IC | | 12/14/2015 19:55 | 1 | | DB-624 0.18(mm) |
| STDTHT 240-210738/21 IC | | 12/14/2015 20:18 | 1 | | DB-624 0.18(mm) |
| STDTHT 240-210738/22 IC | | 12/14/2015 20:40 | 1 | | DB-624 0.18(mm) |

GC/MS VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Instrument ID: A3UX15  Start Date: 01/04/2016 09:30

Analysis Batch Number: 212893  End Date: 01/04/2016 18:08

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| BFB 240-212893/1 | | 01/04/2016 09:30 | 1 | BFB505.D | DB-624 0.18(mm) |
| CCVIS 240-212893/2 | | 01/04/2016 09:55 | 1 | UXC5592.D | DB-624 0.18(mm) |
| LCS 240-212893/5 | | 01/04/2016 10:18 | 1 | UXC5593.D | DB-624 0.18(mm) |
| CCV 240-212893/3 | | 01/04/2016 10:41 | 1 | UXC5594.D | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 11:03 | 1 | | DB-624 0.18(mm) |
| MB 240-212893/8 | | 01/04/2016 11:26 | 1 | UXC5596.D | DB-624 0.18(mm) |
| 240-59482-1 | TB-122815 | 01/04/2016 12:07 | 1 | UXC5597.D | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 12:30 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 12:52 | 66.67 | | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 13:15 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 13:37 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 14:45 | 11.76 | | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 15:07 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 15:30 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 16:15 | 1 | | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 16:38 | 500 | | DB-624 0.18(mm) |
| 240-59545-B-2 MS | | 01/04/2016 17:00 | 500 | UXC5610.D | DB-624 0.18(mm) |
| 240-59545-B-2 MSD | | 01/04/2016 17:23 | 500 | UXC5611.D | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 17:45 | 166.67 | | DB-624 0.18(mm) |
| ZZZZZ | | 01/04/2016 18:08 | 166.67 | | DB-624 0.18(mm) |

# GC/MS VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Batch Number: 212657  Batch Start Date: 12/30/15 14:25  Batch Analyst: Macenczak, Steven

Batch Method: 5030C  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-59482-2 | F-VB-UST-P-7 | 5030C, 8260C | T | 5.26 g | 5 mL | | | | |
| 240-59482-2 MS | F-VB-UST-P-7 | 5030C, 8260C | T | 5.65 g | 5 mL | | | | |
| 240-59482-2 MSD | F-VB-UST-P-7 | 5030C, 8260C | T | 5.47 g | 5 mL | | | | |
| 240-59482-3 | F-VB-UST-P-8 | 5030C, 8260C | T | 5.43 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| | |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8270D

## Semivolatile Organic Compounds (GC/MS) by Method 8270D

FORM II
GC/MS SEMI VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Matrix: Solid  Level: Low

GC Column (1): RXI-5SILMS  ID: 0.45(mm)

| Client Sample ID | Lab Sample ID | 2FP # | PHL # | NBZ # | FBP # | TBP # | TPH # |
|---|---|---|---|---|---|---|---|
| F-VB-UST-P-7 | 240-59482-2 | 81 | 83 | 80 | 77 | 88 | 87 |
| F-VB-UST-P-8 | 240-59482-3 | 83 | 85 | 81 | 77 | 101 | 87 |
| | MB 240-212548/7-A | 81 | 86 | 79 | 78 | 80 | 89 |
| | LCS 240-212548/8-A | 77 | 80 | 82 | 74 | 90 | 81 |
| F-VB-UST-P-7 MS | 240-59482-2 MS | 78 | 80 | 84 | 77 | 101 | 89 |
| F-VB-UST-P-7 MSD | 240-59482-2 MSD | 77 | 79 | 86 | 78 | 90 | 89 |

| | QC LIMITS |
|---|---|
| 2FP = 2-Fluorophenol (Surr) | 24-110 |
| PHL = Phenol-d5 (Surr) | 26-110 |
| NBZ = Nitrobenzene-d5 (Surr) | 20-110 |
| FBP = 2-Fluorobiphenyl (Surr) | 24-110 |
| TBP = 2,4,6-Tribromophenol (Surr) | 10-110 |
| TPH = Terphenyl-d14 (Surr) | 36-110 |

# Column to be used to flag recovery values

FORM II 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51231004.D

Lab ID: LCS 240-212548/8-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 494 | 74 | 35-110 | |
| bis (2-chloroisopropyl) ether | 667 | 493 | 74 | 29-110 | |
| 2,4,5-Trichlorophenol | 667 | 502 | 75 | 25-110 | |
| 2,4,6-Trichlorophenol | 667 | 520 | 78 | 12-110 | |
| 2,4-Dichlorophenol | 667 | 517 | 78 | 39-110 | |
| 2,4-Dimethylphenol | 667 | 505 | 76 | 29-110 | |
| 2,4-Dinitrophenol | 1330 | 563 | 42 | 10-110 | |
| 2,4-Dinitrotoluene | 667 | 582 | 87 | 48-110 | |
| 2,6-Dinitrotoluene | 667 | 568 | 85 | 45-110 | |
| 2-Chloronaphthalene | 667 | 499 | 75 | 32-110 | |
| 2-Chlorophenol | 667 | 524 | 79 | 37-110 | |
| 2-Methylnaphthalene | 667 | 514 | 77 | 36-110 | |
| 2-Methylphenol | 667 | 526 | 79 | 41-110 | |
| 2-Nitroaniline | 667 | 606 | 91 | 45-110 | |
| 2-Nitrophenol | 667 | 577 | 87 | 34-110 | |
| 3,3'-Dichlorobenzidine | 1330 | 1100 | 82 | 28-110 | |
| 3-Nitroaniline | 667 | 536 | 80 | 44-110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 1000 | 75 | 10-110 | |
| 4-Bromophenyl phenyl ether | 667 | 554 | 83 | 39-110 | |
| 4-Chloro-3-methylphenol | 667 | 558 | 84 | 48-110 | |
| 4-Chloroaniline | 667 | 441 | 66 | 30-110 | |
| 4-Chlorophenyl phenyl ether | 667 | 513 | 77 | 40-110 | |
| 4-Nitroaniline | 667 | 630 | 94 | 48-110 | |
| 4-Nitrophenol | 1330 | 1450 | 109 | 28-110 | |
| Acenaphthene | 667 | 501 | 75 | 38-110 | |
| Acenaphthylene | 667 | 491 | 74 | 40-110 | |
| Acetophenone | 667 | 498 | 75 | 40-110 | |
| Anthracene | 667 | 528 | 79 | 48-110 | |
| Atrazine | 1330 | 1330 | 100 | 54-120 | |
| Benzaldehyde | 1330 | 1090 | 82 | 32-110 | |
| Benzo[a]anthracene | 667 | 537 | 80 | 50-110 | |
| Benzo[a]pyrene | 667 | 522 | 78 | 44-110 | |
| Benzo[b]fluoranthene | 667 | 524 | 79 | 43-110 | |
| Benzo[g,h,i]perylene | 667 | 567 | 85 | 51-110 | |
| Benzo[k]fluoranthene | 667 | 525 | 79 | 38-105 | |
| Bis(2-chloroethoxy)methane | 667 | 546 | 82 | 32-110 | |
| Bis(2-chloroethyl)ether | 667 | 501 | 75 | 34-110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 642 | 96 | 50-110 | |
| Butyl benzyl phthalate | 667 | 609 | 91 | 51-110 | |
| Caprolactam | 1330 | 1140 | 85 | 44-114 | |
| Carbazole | 667 | 638 | 96 | 50-110 | |
| Chrysene | 667 | 544 | 82 | 50-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51231004.D

Lab ID: LCS 240-212548/8-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 667 | 581 | 87 | 51-110 | |
| Dibenzofuran | 667 | 503 | 75 | 43-110 | |
| Diethyl phthalate | 667 | 563 | 84 | 52-110 | |
| Dimethyl phthalate | 667 | 550 | 82 | 50-110 | |
| Di-n-butyl phthalate | 667 | 708 | 106 | 51-110 | |
| Di-n-octyl phthalate | 667 | 571 | 86 | 48-110 | |
| Fluoranthene | 667 | 589 | 88 | 51-110 | |
| Fluorene | 667 | 465 | 70 | 46-110 | |
| Hexachlorobenzene | 667 | 523 | 79 | 43-110 | |
| Hexachlorobutadiene | 667 | 515 | 77 | 29-110 | |
| Hexachlorocyclopentadiene | 667 | 399 | 60 | 12-110 | |
| Hexachloroethane | 667 | 520 | 78 | 30-110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 574 | 86 | 50-110 | |
| Isophorone | 667 | 555 | 83 | 36-110 | |
| N-Nitrosodi-n-propylamine | 667 | 544 | 82 | 38-110 | |
| N-Nitrosodiphenylamine | 1330 | 1100 | 82 | 46-110 | |
| Naphthalene | 667 | 501 | 75 | 36-110 | |
| Nitrobenzene | 667 | 545 | 82 | 32-110 | |
| Pentachlorophenol | 1330 | 1110 | 83 | 10-110 | |
| Phenanthrene | 667 | 530 | 79 | 49-110 | |
| Phenol | 667 | 505 | 76 | 38-110 | |
| Pyrene | 667 | 574 | 86 | 49-110 | |
| 3 & 4 Methylphenol | 667 | 516 | 77 | 40-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51231009.D

Lab ID: 240-59482-2 MS Client ID: F-VB-UST-P-7 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 826 | 4.3 U | 631 | 76 | 32-110 | |
| bis (2-chloroisopropyl) ether | 826 | 12 U | 607 | 73 | 11-110 | |
| 2,4,5-Trichlorophenol | 826 | 31 U | 627 | 76 | 10-117 | |
| 2,4,6-Trichlorophenol | 826 | 11 U | 670 | 81 | 10-110 | |
| 2,4-Dichlorophenol | 826 | 24 U | 640 | 77 | 10-110 | |
| 2,4-Dimethylphenol | 826 | 24 U | 717 | 87 | 10-110 | |
| 2,4-Dinitrophenol | 1650 | 26 U | 1550 | 94 | 10-110 | |
| 2,4-Dinitrotoluene | 826 | 21 U | 788 | 95 | 32-110 | |
| 2,6-Dinitrotoluene | 826 | 26 U | 752 | 91 | 35-110 | |
| 2-Chloronaphthalene | 826 | 0.55 U | 636 | 77 | 28-110 | |
| 2-Chlorophenol | 826 | 10 U | 631 | 76 | 10-110 | |
| 2-Methylnaphthalene | 826 | 0.61 U | 629 | 76 | 10-133 | |
| 2-Methylphenol | 826 | 13 U | 623 | 75 | 24-110 | |
| 2-Nitroaniline | 826 | 11 U | 808 | 98 | 39-110 | |
| 2-Nitrophenol | 826 | 10 U | 710 | 86 | 10-110 | |
| 3,3'-Dichlorobenzidine | 1650 | 22 U | 1320 | 80 | 10-110 | |
| 3-Nitroaniline | 826 | 20 U | 490 | 59 | 10-110 | |
| 4,6-Dinitro-2-methylphenol | 1650 | 11 U | 1620 | 98 | 10-110 | |
| 4-Bromophenyl phenyl ether | 826 | 16 U | 767 | 93 | 33-110 | |
| 4-Chloro-3-methylphenol | 826 | 26 U | 727 | 88 | 25-110 | |
| 4-Chloroaniline | 826 | 21 U | 396 | 48 | 10-110 | |
| 4-Chlorophenyl phenyl ether | 826 | 16 U | 672 | 81 | 32-110 | |
| 4-Nitroaniline | 826 | 32 U | 664 | 80 | 10-110 | |
| 4-Nitrophenol | 1650 | 21 U | 2010 | 122 | 10-113 | F1 |
| Acenaphthene | 826 | 0.93 U | 648 | 78 | 22-110 | |
| Acenaphthylene | 826 | 0.43 U | 654 | 79 | 24-110 | |
| Acetophenone | 826 | 11 U | 652 | 79 | 31-110 | |
| Anthracene | 826 | 0.95 U | 711 | 86 | 20-110 | |
| Atrazine | 1650 | 11 U | 1730 | 105 | 45-118 | |
| Benzaldehyde | 1650 | 15 U | 1330 | 81 | 23-110 | |
| Benzo[a]anthracene | 826 | 0.77 U | 711 | 86 | 10-122 | |
| Benzo[a]pyrene | 826 | 0.78 U | 713 | 86 | 10-110 | |
| Benzo[b]fluoranthene | 826 | 0.72 U | 746 | 90 | 12-118 | |
| Benzo[g,h,i]perylene | 826 | 0.43 U | 819 | 99 | 10-117 | |
| Benzo[k]fluoranthene | 826 | 0.83 U | 712 | 86 | 10-121 | |
| Bis(2-chloroethoxy)methane | 826 | 27 U | 681 | 82 | 26-110 | |
| Bis(2-chloroethyl)ether | 826 | 2.4 U | 697 | 84 | 21-110 | |
| Bis(2-ethylhexyl) phthalate | 826 | 160 | 861 | 85 | 40-110 | |
| Butyl benzyl phthalate | 826 | 12 U | 822 | 100 | 44-110 | |
| Caprolactam | 1650 | 45 U | 1490 | 90 | 10-134 | |
| Carbazole | 826 | 33 U | 763 | 92 | 34-110 | |
| Chrysene | 826 | 1.3 U | 717 | 87 | 10-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51231009.D

Lab ID: 240-59482-2 MS Client ID: F-VB-UST-P-7 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 826 | 0.81 U | 844 | 102 | 14-113 | |
| Dibenzofuran | 826 | 0.81 U | 669 | 81 | 29-110 | |
| Diethyl phthalate | 826 | 20 U | 779 | 94 | 42-110 | |
| Dimethyl phthalate | 826 | 21 U | 741 | 90 | 41-110 | |
| Di-n-butyl phthalate | 826 | 30 J | 937 | 110 | 43-110 | |
| Di-n-octyl phthalate | 826 | 9.7 U | 794 | 96 | 24-119 | |
| Fluoranthene | 826 | 9.3 | 781 | 93 | 10-110 | |
| Fluorene | 826 | 0.65 U | 631 | 76 | 23-110 | |
| Hexachlorobenzene | 826 | 2.6 U | 709 | 86 | 34-110 | |
| Hexachlorobutadiene | 826 | 6.9 U | 629 | 76 | 25-110 | |
| Hexachlorocyclopentadiene | 826 | 9.9 U | 333 J | 40 | 10-110 | |
| Hexachloroethane | 826 | 11 U | 617 | 75 | 12-110 | |
| Indeno[1,2,3-cd]pyrene | 826 | 0.43 U | 811 | 98 | 10-114 | |
| Isophorone | 826 | 21 J | 676 | 79 | 29-110 | |
| N-Nitrosodi-n-propylamine | 826 | 7.7 U | 674 | 82 | 26-110 | |
| N-Nitrosodiphenylamine | 1650 | 26 U | 1440 | 87 | 22-110 | |
| Naphthalene | 826 | 4.3 J | 612 | 74 | 10-111 | |
| Nitrobenzene | 826 | 2.7 U | 676 | 82 | 23-110 | |
| Pentachlorophenol | 1650 | 11 U | 1520 | 92 | 10-110 | |
| Phenanthrene | 826 | 9.3 | 715 | 85 | 10-166 | |
| Phenol | 826 | 8.9 U | 601 | 73 | 17-110 | |
| Pyrene | 826 | 7.9 J | 773 | 93 | 10-147 | |
| 3 & 4 Methylphenol | 826 | 24 U | 613 | 74 | 25-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51231010.D

Lab ID: 240-59482-2 MSD Client ID: F-VB-UST-P-7 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 816 | 636 | 78 | 1 | 32 | 32-110 | |
| bis (2-chloroisopropyl) ether | 816 | 617 | 76 | 2 | 42 | 11-110 | |
| 2,4,5-Trichlorophenol | 816 | 643 | 79 | 3 | 99 | 10-117 | |
| 2,4,6-Trichlorophenol | 816 | 636 | 78 | 5 | 38 | 10-110 | |
| 2,4-Dichlorophenol | 816 | 646 | 79 | 1 | 34 | 10-110 | |
| 2,4-Dimethylphenol | 816 | 684 | 84 | 5 | 31 | 10-110 | |
| 2,4-Dinitrophenol | 1630 | 1450 | 89 | 7 | 99 | 10-110 | |
| 2,4-Dinitrotoluene | 816 | 801 | 98 | 2 | 30 | 32-110 | |
| 2,6-Dinitrotoluene | 816 | 742 | 91 | 1 | 30 | 35-110 | |
| 2-Chloronaphthalene | 816 | 636 | 78 | 0 | 30 | 28-110 | |
| 2-Chlorophenol | 816 | 640 | 78 | 1 | 47 | 10-110 | |
| 2-Methylnaphthalene | 816 | 641 | 78 | 2 | 42 | 10-133 | |
| 2-Methylphenol | 816 | 638 | 78 | 2 | 51 | 24-110 | |
| 2-Nitroaniline | 816 | 791 | 97 | 2 | 31 | 39-110 | |
| 2-Nitrophenol | 816 | 752 | 92 | 6 | 49 | 10-110 | |
| 3,3'-Dichlorobenzidine | 1630 | 1470 | 90 | 11 | 56 | 10-110 | |
| 3-Nitroaniline | 816 | 589 | 72 | 18 | 30 | 10-110 | |
| 4,6-Dinitro-2-methylphenol | 1630 | 1590 | 97 | 2 | 55 | 10-110 | |
| 4-Bromophenyl phenyl ether | 816 | 736 | 90 | 4 | 30 | 33-110 | |
| 4-Chloro-3-methylphenol | 816 | 726 | 89 | 0 | 54 | 25-110 | |
| 4-Chloroaniline | 816 | 378 | 46 | 5 | 36 | 10-110 | |
| 4-Chlorophenyl phenyl ether | 816 | 665 | 81 | 1 | 30 | 32-110 | |
| 4-Nitroaniline | 816 | 799 | 98 | 18 | 48 | 10-110 | |
| 4-Nitrophenol | 1630 | 1740 | 107 | 14 | 49 | 10-113 | |
| Acenaphthene | 816 | 636 | 78 | 2 | 99 | 22-110 | |
| Acenaphthylene | 816 | 646 | 79 | 1 | 99 | 24-110 | |
| Acetophenone | 816 | 648 | 79 | 1 | 43 | 31-110 | |
| Anthracene | 816 | 704 | 86 | 1 | 99 | 20-110 | |
| Atrazine | 1630 | 1740 | 107 | 1 | 30 | 45-118 | |
| Benzaldehyde | 1630 | 1270 | 78 | 4 | 42 | 23-110 | |
| Benzo[a]anthracene | 816 | 731 | 90 | 3 | 99 | 10-122 | |
| Benzo[a]pyrene | 816 | 714 | 87 | 0 | 99 | 10-110 | |
| Benzo[b]fluoranthene | 816 | 738 | 90 | 1 | 99 | 12-118 | |
| Benzo[g,h,i]perylene | 816 | 783 | 96 | 4 | 99 | 10-117 | |
| Benzo[k]fluoranthene | 816 | 716 | 88 | 1 | 99 | 10-121 | |
| Bis(2-chloroethoxy)methane | 816 | 683 | 84 | 0 | 37 | 26-110 | |
| Bis(2-chloroethyl)ether | 816 | 637 | 78 | 9 | 55 | 21-110 | |
| Bis(2-ethylhexyl) phthalate | 816 | 938 | 95 | 9 | 30 | 40-110 | |
| Butyl benzyl phthalate | 816 | 832 | 102 | 1 | 30 | 44-110 | |
| Caprolactam | 1630 | 1500 | 92 | 1 | 32 | 10-134 | |
| Carbazole | 816 | 828 | 101 | 8 | 30 | 34-110 | |
| Chrysene | 816 | 734 | 90 | 2 | 99 | 10-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51231010.D

Lab ID: 240-59482-2 MSD Client ID: F-VB-UST-P-7 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 816 | 805 | 99 | 5 | 99 | 14-113 | |
| Dibenzofuran | 816 | 656 | 80 | 2 | 30 | 29-110 | |
| Diethyl phthalate | 816 | 764 | 94 | 2 | 30 | 42-110 | |
| Dimethyl phthalate | 816 | 730 | 89 | 1 | 30 | 41-110 | |
| Di-n-butyl phthalate | 816 | 939 | 111 | 0 | 30 | 43-110 | F1 |
| Di-n-octyl phthalate | 816 | 788 | 97 | 1 | 30 | 24-119 | |
| Fluoranthene | 816 | 787 | 95 | 1 | 99 | 10-110 | |
| Fluorene | 816 | 619 | 76 | 2 | 99 | 23-110 | |
| Hexachlorobenzene | 816 | 682 | 83 | 4 | 30 | 34-110 | |
| Hexachlorobutadiene | 816 | 683 | 84 | 8 | 34 | 25-110 | |
| Hexachlorocyclopentadiene | 816 | 302 J | 37 | 10 | 79 | 10-110 | |
| Hexachloroethane | 816 | 646 | 79 | 5 | 50 | 12-110 | |
| Indeno[1,2,3-cd]pyrene | 816 | 782 | 96 | 4 | 99 | 10-114 | |
| Isophorone | 816 | 691 | 82 | 2 | 38 | 29-110 | |
| N-Nitrosodi-n-propylamine | 816 | 684 | 84 | 1 | 42 | 26-110 | |
| N-Nitrosodiphenylamine | 1630 | 1420 | 87 | 2 | 30 | 22-110 | |
| Naphthalene | 816 | 623 | 76 | 2 | 99 | 10-111 | |
| Nitrobenzene | 816 | 680 | 83 | 1 | 41 | 23-110 | |
| Pentachlorophenol | 1630 | 1280 | 79 | 17 | 50 | 10-110 | |
| Phenanthrene | 816 | 706 | 85 | 1 | 99 | 10-166 | |
| Phenol | 816 | 623 | 76 | 4 | 53 | 17-110 | |
| Pyrene | 816 | 766 | 93 | 1 | 99 | 10-147 | |
| 3 & 4 Methylphenol | 816 | 618 | 76 | 1 | 50 | 25-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM IV
GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab File ID: 51231003.D  Lab Sample ID: MB 240-212548/7-A

Matrix: Solid  Date Extracted: 12/30/2015 07:05

Instrument ID: A4HP9  Date Analyzed: 12/31/2015 10:02

Level:(Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-212548/8-A | 51231004.D | 12/31/2015 10:53 |
| F-VB-UST-P-8 | 240-59482-3 | 51231007.D | 12/31/2015 12:08 |
| F-VB-UST-P-7 | 240-59482-2 | 51231008.D | 12/31/2015 12:33 |
| F-VB-UST-P-7 MS | 240-59482-2 MS | 51231009.D | 12/31/2015 12:58 |
| F-VB-UST-P-7 MSD | 240-59482-2 MSD | 51231010.D | 12/31/2015 13:23 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab File ID: 51222001.D DFTPP Injection Date: 12/22/2015

Instrument ID: A4HP9 DFTPP Injection Time: 12:26

Analysis Batch No.: 211933

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 56.1 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0) 1 |
| 69 | Mass 69 relative abundance | 59.9 | |
| 70 | Less than 2.0% of mass 69 | 0.3 | (0.4) 1 |
| 127 | 25.0 - 75.0% of mass 198 | 63.4 | |
| 197 | Less than 1.0% of mass 198 | 0.2 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.6 | |
| 275 | 10.0- 30.0% of mass 198 | 21.7 | |
| 365 | Greater than 0.75% of mass 198 | 2.8 | |
| 441 | Present, but less than mass 443 | 10.0 | |
| 442 | 40.0 - 110.0% of mass 198 | 61.9 | |
| 443 | 15.0 - 24.0% of mass 442 | 12.0 | (19.3) 2 |

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD5 240-211933/2 | 51222002.D | 12/22/2015 | 12:48 |
| | STD4 240-211933/3 | 51222003.D | 12/22/2015 | 13:13 |
| | STD3 240-211933/4 | 51222004.D | 12/22/2015 | 13:39 |
| | STD2 240-211933/5 | 51222005.D | 12/22/2015 | 14:04 |
| | STD1 240-211933/6 | 51222006.D | 12/22/2015 | 14:30 |
| | STD9 240-211933/7 | 51222007.D | 12/22/2015 | 14:56 |
| | STD8 240-211933/8 | 51222008.D | 12/22/2015 | 15:21 |
| | STD7 240-211933/9 | 51222009.D | 12/22/2015 | 15:47 |
| | STD6 240-211933/10 | 51222010.D | 12/22/2015 | 16:12 |
| | ICV 240-211933/11 | 51222011.D | 12/22/2015 | 16:38 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab File ID: 51231001.D  DFTPP Injection Date: 12/31/2015

Instrument ID: A4HP9  DFTPP Injection Time: 09:15

Analysis Batch No.: 212723

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 62.8 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0) 1 |
| 69 | Mass 69 relative abundance | 65.2 | |
| 70 | Less than 2.0% of mass 69 | 0.3 | (0.5) 1 |
| 127 | 25.0 - 75.0% of mass 198 | 71.9 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.9 | |
| 275 | 10.0- 30.0% of mass 198 | 23.7 | |
| 365 | Greater than 0.75% of mass 198 | 3.7 | |
| 441 | Present, but less than mass 443 | 13.8 | |
| 442 | 40.0 - 110.0% of mass 198 | 87.2 | |
| 443 | 15.0 - 24.0% of mass 442 | 16.6 | (19.1) 2 |

1-Value is % mass 69  2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCV 240-212723/2 | 51231002.D | 12/31/2015 | 09:36 |
| | MB 240-212548/7-A | 51231003.D | 12/31/2015 | 10:02 |
| | LCS 240-212548/8-A | 51231004.D | 12/31/2015 | 10:53 |
| F-VB-UST-P-8 | 240-59482-3 | 51231007.D | 12/31/2015 | 12:08 |
| F-VB-UST-P-7 | 240-59482-2 | 51231008.D | 12/31/2015 | 12:33 |
| F-VB-UST-P-7 MS | 240-59482-2 MS | 51231009.D | 12/31/2015 | 12:58 |
| F-VB-UST-P-7 MSD | 240-59482-2 MSD | 51231010.D | 12/31/2015 | 13:23 |

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Sample No.: STD6 240-211933/10 Date Analyzed: 12/22/2015 16:12

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm)

Lab File ID (Standard): 51222010.D Heated Purge: (Y/N) N

Calibration ID: 32225

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 239791 | 6.57 | 975433 | 7.68 | 511807 | 9.19 |
| UPPER LIMIT | | 479582 | 7.07 | 1950866 | 8.18 | 1023614 | 9.69 |
| LOWER LIMIT | | 119896 | 6.07 | 487717 | 7.18 | 255904 | 8.69 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-211933/11 | | 264514 | 6.57 | 1093427 | 7.68 | 567798 | 9.19 |
| CCV 240-212723/2 CCVIS | | 218841 | 6.47 | 880111 | 7.59 | 457488 | 9.09 |

DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Sample No.: STD6 240-211933/10 Date Analyzed: 12/22/2015 16:12

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm)

Lab File ID (Standard): 51222010.D Heated Purge: (Y/N) N

Calibration ID: 32225

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 747104 | 10.46 | 847010 | 12.81 | 728741 | 14.76 |
| UPPER LIMIT | | 1494208 | 10.96 | 1694020 | 13.31 | 1457482 | 15.26 |
| LOWER LIMIT | | 373552 | 9.96 | 423505 | 12.31 | 364371 | 14.26 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-211933/11 | | 847727 | 10.46 | 949631 | 12.81 | 828488 | 14.76 |
| CCV 240-212723/2 CCVIS | | 698427 | 10.36 | 724352 | 12.67 | 651306 | 14.52 |

PHN = Phenanthrene-d10
CRY = Chrysene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Sample No.: CCV 240-212723/2 Date Analyzed: 12/31/2015 09:36

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm)

Lab File ID (Standard): 51231002.D Heated Purge: (Y/N) N

Calibration ID: 32229

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 218841 | 6.47 | 880111 | 7.59 | 457488 | 9.09 |
| UPPER LIMIT | | 437682 | 6.97 | 1760222 | 8.09 | 914976 | 9.59 |
| LOWER LIMIT | | 109421 | 5.97 | 440056 | 7.09 | 228744 | 8.59 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-212548/7-A | | 163289 | 6.47 | 698467 | 7.58 | 371239 | 9.09 |
| LCS 240-212548/8-A | | 191356 | 6.47 | 779708 | 7.58 | 422992 | 9.09 |
| 240-59482-3 | F-VB-UST-P-8 | 168300 | 6.47 | 688973 | 7.58 | 370596 | 9.08 |
| 240-59482-2 | F-VB-UST-P-7 | 174672 | 6.47 | 714524 | 7.58 | 382972 | 9.09 |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 169885 | 6.47 | 690559 | 7.58 | 362728 | 9.09 |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 179051 | 6.47 | 726872 | 7.58 | 390035 | 9.09 |

DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Sample No.: CCV 240-212723/2  Date Analyzed: 12/31/2015 09:36

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)

Lab File ID (Standard): 51231002.D  Heated Purge: (Y/N) N

Calibration ID: 32229

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 698427 | 10.36 | 724352 | 12.67 | 651306 | 14.52 |
| UPPER LIMIT | | 1396854 | 10.86 | 1448704 | 13.17 | 1302612 | 15.02 |
| LOWER LIMIT | | 349214 | 9.86 | 362176 | 12.17 | 325653 | 14.02 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-212548/7-A | | 583813 | 10.36 | 592138 | 12.66 | 528025 | 14.52 |
| LCS 240-212548/8-A | | 631254 | 10.36 | 681934 | 12.67 | 618178 | 14.52 |
| 240-59482-3 | F-VB-UST-P-8 | 574436 | 10.36 | 606579 | 12.66 | 544844 | 14.52 |
| 240-59482-2 | F-VB-UST-P-7 | 608076 | 10.36 | 617593 | 12.67 | 544288 | 14.52 |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 552277 | 10.36 | 586705 | 12.67 | 525748 | 14.52 |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 597920 | 10.36 | 637615 | 12.67 | 575735 | 14.52 |

PHN = Phenanthrene-d10
CRY = Chrysene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 | Lab Sample ID: 240-59482-2

Matrix: Solid | Lab File ID: 51231008.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:00

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30.06(g) | Date Analyzed: 12/31/2015 12:33

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.3 | U | 61 | 4.3 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 12 | U | 120 | 12 |
| 95-95-4 | 2,4,5-Trichlorophenol | 31 | U | 180 | 31 |
| 88-06-2 | 2,4,6-Trichlorophenol | 11 | U | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 26 | U | 400 | 26 |
| 121-14-2 | 2,4-Dinitrotoluene | 21 | U | 240 | 21 |
| 606-20-2 | 2,6-Dinitrotoluene | 26 | U | 240 | 26 |
| 91-58-7 | 2-Chloronaphthalene | 0.55 | U | 61 | 0.55 |
| 95-57-8 | 2-Chlorophenol | 10 | U | 61 | 10 |
| 91-57-6 | 2-Methylnaphthalene | 0.61 | U | 8.2 | 0.61 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 10 | U | 61 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 22 | U | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 20 | U | 240 | 20 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 16 | U | 61 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 26 | U | 180 | 26 |
| 106-47-8 | 4-Chloroaniline | 21 | U | 180 | 21 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 16 | U | 61 | 16 |
| 100-01-6 | 4-Nitroaniline | 32 | U | 240 | 32 |
| 100-02-7 | 4-Nitrophenol | 21 | U F1 | 400 | 21 |
| 83-32-9 | Acenaphthene | 0.93 | U | 8.2 | 0.93 |
| 208-96-8 | Acenaphthylene | 0.43 | U | 8.2 | 0.43 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 0.95 | U | 8.2 | 0.95 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 15 | U | 120 | 15 |
| 56-55-3 | Benzo[a]anthracene | 0.77 | U | 8.2 | 0.77 |
| 50-32-8 | Benzo[a]pyrene | 0.78 | U | 8.2 | 0.78 |
| 205-99-2 | Benzo[b]fluoranthene | 0.72 | U | 8.2 | 0.72 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.43 | U | 8.2 | 0.43 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 | Lab Sample ID: 240-59482-2

Matrix: Solid | Lab File ID: 51231008.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:00

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30.06(g) | Date Analyzed: 12/31/2015 12:33

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.83 | U | 8.2 | 0.83 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 27 | U | 120 | 27 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 160 | B | 86 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 86 | 12 |
| 105-60-2 | Caprolactam | 45 | U | 400 | 45 |
| 86-74-8 | Carbazole | 33 | U | 61 | 33 |
| 218-01-9 | Chrysene | 1.3 | U | 8.2 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.81 | U | 8.2 | 0.81 |
| 132-64-9 | Dibenzofuran | 0.81 | U | 61 | 0.81 |
| 84-66-2 | Diethyl phthalate | 20 | U | 86 | 20 |
| 131-11-3 | Dimethyl phthalate | 21 | U | 86 | 21 |
| 84-74-2 | Di-n-butyl phthalate | 30 | J B F1 | 86 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.7 | U | 86 | 9.7 |
| 206-44-0 | Fluoranthene | 9.3 | | 8.2 | 0.67 |
| 86-73-7 | Fluorene | 0.65 | U | 8.2 | 0.65 |
| 118-74-1 | Hexachlorobenzene | 2.6 | U | 8.2 | 2.6 |
| 87-68-3 | Hexachlorobutadiene | 6.9 | U | 61 | 6.9 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.9 | U | 400 | 9.9 |
| 67-72-1 | Hexachloroethane | 11 | U | 61 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.43 | U | 8.2 | 0.43 |
| 78-59-1 | Isophorone | 21 | J B | 61 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.7 | U | 61 | 7.7 |
| 86-30-6 | N-Nitrosodiphenylamine | 26 | U | 61 | 26 |
| 91-20-3 | Naphthalene | 4.3 | J | 8.2 | 1.0 |
| 98-95-3 | Nitrobenzene | 2.7 | U | 120 | 2.7 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 9.3 | | 8.2 | 0.89 |
| 108-95-2 | Phenol | 8.9 | U | 61 | 8.9 |
| 129-00-0 | Pyrene | 7.9 | J | 8.2 | 0.54 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 490 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 | Lab Sample ID: 240-59482-2

Matrix: Solid | Lab File ID: 51231008.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:00

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30.06(g) | Date Analyzed: 12/31/2015 12:33

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 87 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 83 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 80 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 81 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 77 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 88 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-8 | Lab Sample ID: 240-59482-3

Matrix: Solid | Lab File ID: 51231007.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:30

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30.35(g) | Date Analyzed: 12/31/2015 12:08

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 4.1 | U | 59 | 4.1 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 11 | U | 120 | 11 |
| 95-95-4 | 2,4,5-Trichlorophenol | 29 | U | 180 | 29 |
| 88-06-2 | 2,4,6-Trichlorophenol | 10 | U | 180 | 10 |
| 120-83-2 | 2,4-Dichlorophenol | 24 | U | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 24 | U | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 25 | U | 390 | 25 |
| 121-14-2 | 2,4-Dinitrotoluene | 20 | U | 240 | 20 |
| 606-20-2 | 2,6-Dinitrotoluene | 25 | U | 240 | 25 |
| 91-58-7 | 2-Chloronaphthalene | 0.53 | U | 59 | 0.53 |
| 95-57-8 | 2-Chlorophenol | 9.7 | U | 59 | 9.7 |
| 91-57-6 | 2-Methylnaphthalene | 53 | | 7.9 | 0.59 |
| 95-48-7 | 2-Methylphenol | 13 | U | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 11 | U | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 9.8 | U | 59 | 9.8 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 21 | U | 120 | 21 |
| 99-09-2 | 3-Nitroaniline | 19 | U | 240 | 19 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 11 | U | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 15 | U | 59 | 15 |
| 59-50-7 | 4-Chloro-3-methylphenol | 25 | U | 180 | 25 |
| 106-47-8 | 4-Chloroaniline | 20 | U | 180 | 20 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 15 | U | 59 | 15 |
| 100-01-6 | 4-Nitroaniline | 31 | U | 240 | 31 |
| 100-02-7 | 4-Nitrophenol | 20 | U | 390 | 20 |
| 83-32-9 | Acenaphthene | 63 | | 7.9 | 0.90 |
| 208-96-8 | Acenaphthylene | 0.41 | U | 7.9 | 0.41 |
| 98-86-2 | Acetophenone | 11 | U | 120 | 11 |
| 120-12-7 | Anthracene | 97 | | 7.9 | 0.92 |
| 1912-24-9 | Atrazine | 11 | U | 240 | 11 |
| 100-52-7 | Benzaldehyde | 14 | U | 120 | 14 |
| 56-55-3 | Benzo[a]anthracene | 140 | | 7.9 | 0.74 |
| 50-32-8 | Benzo[a]pyrene | 100 | | 7.9 | 0.75 |
| 205-99-2 | Benzo[b]fluoranthene | 180 | | 7.9 | 0.70 |
| 191-24-2 | Benzo[g,h,i]perylene | 51 | | 7.9 | 0.41 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-8 | Lab Sample ID: 240-59482-3

Matrix: Solid | Lab File ID: 51231007.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:30

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30.35(g) | Date Analyzed: 12/31/2015 12:08

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 47 | | 7.9 | 0.80 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 26 | U | 120 | 26 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4 | U | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 82 | B | 82 | 22 |
| 85-68-7 | Butyl benzyl phthalate | 12 | U | 82 | 12 |
| 105-60-2 | Caprolactam | 44 | U | 390 | 44 |
| 86-74-8 | Carbazole | 32 | U | 59 | 32 |
| 218-01-9 | Chrysene | 94 | | 7.9 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 16 | | 7.9 | 0.78 |
| 132-64-9 | Dibenzofuran | 51 | J | 59 | 0.78 |
| 84-66-2 | Diethyl phthalate | 19 | U | 82 | 19 |
| 131-11-3 | Dimethyl phthalate | 20 | U | 82 | 20 |
| 84-74-2 | Di-n-butyl phthalate | 23 | J B | 82 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 9.3 | U | 82 | 9.3 |
| 206-44-0 | Fluoranthene | 490 | | 7.9 | 0.65 |
| 86-73-7 | Fluorene | 88 | | 7.9 | 0.62 |
| 118-74-1 | Hexachlorobenzene | 2.5 | U | 7.9 | 2.5 |
| 87-68-3 | Hexachlorobutadiene | 6.6 | U | 59 | 6.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 9.5 | U | 390 | 9.5 |
| 67-72-1 | Hexachloroethane | 11 | U | 59 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 45 | | 7.9 | 0.41 |
| 78-59-1 | Isophorone | 20 | J B | 59 | 15 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 7.4 | U | 59 | 7.4 |
| 86-30-6 | N-Nitrosodiphenylamine | 25 | U | 59 | 25 |
| 91-20-3 | Naphthalene | 14 | | 7.9 | 0.97 |
| 98-95-3 | Nitrobenzene | 2.6 | U | 120 | 2.6 |
| 87-86-5 | Pentachlorophenol | 11 | U | 180 | 11 |
| 85-01-8 | Phenanthrene | 570 | | 7.9 | 0.86 |
| 108-95-2 | Phenol | 8.6 | U | 59 | 8.6 |
| 129-00-0 | Pyrene | 370 | | 7.9 | 0.52 |
| 15831-10-4 | 3 & 4 Methylphenol | 24 | U | 470 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-8 | Lab Sample ID: 240-59482-3

Matrix: Solid | Lab File ID: 51231007.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:30

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30.35(g) | Date Analyzed: 12/31/2015 12:08

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 87 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 85 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 81 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 83 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 77 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 101 | | 10-110 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-59482-1  Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48  Calibration End Date: 12/22/2015 20:02  Calibration ID: 32225

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-211933/6 | 51222006.D |
| Level 2 | STD2 240-211933/5 | 51222005.D |
| Level 3 | STD3 240-211933/4 | 51222004.D |
| Level 4 | STD4 240-211933/3 | 51222003.D |
| Level 5 | STD5 240-211933/2 | 51222002.D |
| Level 6 | STD6 240-211933/10 | 51222010.D |
| Level 7 | STD7 240-211933/9 | 51222009.D |
| Level 8 | STD8 240-211933/8 | 51222008.D |
| Level 9 | STD9 240-211933/7 | 51222007.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,4-Dioxane | 0.7493 | 0.6839<br>0.6809 | 0.7228<br>0.7301 | 0.6844<br>0.7284 | 0.6965 | Ave | | 0.7095 | | | | 3.7 | | 20.0 | | | |
| N-Nitrosodimethylamine | 1.1344 | 1.0072<br>1.0301 | 1.0045<br>1.0825 | 1.0386<br>1.0934 | 0.9948 | Ave | | 1.0482 | | | | 4.8 | | 20.0 | | | |
| Pyridine | 2.0061 | 1.8287<br>1.8457 | 1.9380<br>1.9040 | 1.7810<br>1.9202 | 1.7658 | Ave | | 1.8737 | | | | 4.4 | | 20.0 | | | |
| Benzaldehyde | 1.2874 | 1.5081<br>1.2814 | 1.4998<br>1.2478 | 1.5369<br>1.2039 | 1.4064 | Ave | | 1.3715 | | | 0.0100 | 9.6 | | 20.0 | | | |
| Phenol | 2.0257 | 2.0966<br>1.9251 | 2.1688<br>1.9217 | 2.1149<br>1.9052 | 2.0095 | Ave | | 2.0209 | | | 0.8000 | 4.9 | | 20.0 | | | |
| Aniline | 2.5558 | 2.5916<br>2.3743 | 2.5576<br>2.4056 | 2.5746<br>2.4173 | 2.4724 | Ave | | 2.4937 | | | | 3.5 | | 20.0 | | | |
| Bis(2-chloroethyl)ether | 1.7274 | 1.6519<br>1.6091 | 1.7003<br>1.6112 | 1.7457<br>1.6210 | 1.7057 | Ave | | 1.6715 | | | 0.7000 | 3.3 | | 20.0 | | | |
| 2-Chlorophenol | 1.4519 | 1.4132<br>1.3362 | 1.4989<br>1.3531 | 1.4735<br>1.3411 | 1.4423 | Ave | | 1.4138 | | | 0.8000 | 4.5 | | 20.0 | | | |
| n-Decane | 1.8835 | 1.8319<br>1.6942 | 1.9312<br>1.7005 | 1.9492<br>1.6748 | 1.8157 | Ave | | 1.8101 | | | | 6.0 | | 20.0 | | | |
| 1,3-Dichlorobenzene | 1.5274 | 1.6090<br>1.4115 | 1.6418<br>1.4277 | 1.5661<br>1.4065 | 1.5317 | Ave | | 1.5152 | | | | 6.0 | | 20.0 | | | |
| 1,4-Dichlorobenzene | 1.5125 | 1.5339<br>1.4264 | 1.5981<br>1.4343 | 1.6019<br>1.4281 | 1.5187 | Ave | | 1.5067 | | | | 4.8 | | 20.0 | | | |
| Benzyl alcohol | 1.0913 | 1.1466<br>1.0345 | 1.0998<br>1.0239 | 1.1180<br>1.0413 | 1.0742 | Ave | | 1.0787 | | | | 4.0 | | 20.0 | | | |
| 1,2-Dichlorobenzene | 1.4687 | 1.5441<br>1.3505 | 1.5404<br>1.3691 | 1.5544<br>1.3372 | 1.4774 | Ave | | 1.4552 | | | | 6.3 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48 Calibration End Date: 12/22/2015 20:02 Calibration ID: 32225

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2-Methylphenol | 1.4961 | 1.4796<br>1.4050 | 1.5571<br>1.4137 | 1.5391<br>1.4035 | 1.4702 | Ave | | 1.4705 | | | 0.7000 | 4.1 | | 20.0 | | | |
| bis (2-chloroisopropyl) ether | 2.1439 | 2.1725<br>1.9981 | 2.1847<br>1.9903 | 2.1466<br>1.9760 | 2.0806 | Ave | | 2.0866 | | | | 4.2 | | 20.0 | | | |
| Indene | 2.5272 | 2.7747<br>2.2980 | 2.7981<br>2.2839 | 2.6994<br>2.2504 | 2.5810 | Ave | | 2.5266 | | | | 8.9 | | 20.0 | | | |
| 3 & 4 Methylphenol | 1.5751 | 1.4984<br>1.4738 | 1.6094<br>1.4656 | 1.6010<br>1.4612 | 1.5658 | Ave | | 1.5313 | | | | 4.1 | | 20.0 | | | |
| N-Nitrosodi-n-propylamine | 1.2195 | 1.2913<br>1.1501 | 1.2001<br>1.1297 | 1.2667<br>1.1570 | 1.2461 | Ave | | 1.2076 | | | 0.5000 | 4.9 | | 20.0 | | | |
| Acetophenone | 2.1849 | 2.3275<br>2.0753 | 2.2584<br>2.0366 | 2.2796<br>2.0407 | 2.2198 | Ave | | 2.1779 | | | 0.0100 | 5.2 | | 20.0 | | | |
| Hexachloroethane | 0.6306 | 0.6721<br>0.5942 | 0.6774<br>0.6039 | 0.6590<br>0.5952 | 0.6353 | Ave | | 0.6335 | | | 0.3000 | 5.3 | | 20.0 | | | |
| Nitrobenzene | 0.4704 | 0.4753<br>0.4483 | 0.4698<br>0.4388 | 0.4853<br>0.4354 | 0.4638 | Ave | | 0.4609 | | | 0.2000 | 3.9 | | 20.0 | | | |
| Isophorone | 0.8350 | 0.8495<br>0.7955 | 0.8419<br>0.7695 | 0.8678<br>0.7638 | 0.8374 | Ave | | 0.8201 | | | 0.4000 | 4.7 | | 20.0 | | | |
| 2,4-Dimethylphenol | 0.3840 | 0.4049<br>0.3678 | 0.4119<br>0.3520 | 0.4104<br>0.3447 | 0.3997 | Ave | | 0.3844 | | | 0.2000 | 6.9 | | 20.0 | | | |
| 2-Nitrophenol | 0.1685 | 0.1584<br>0.1590 | 0.1611<br>0.1520 | 0.1682<br>0.1541 | 0.1678 | Ave | | 0.1611 | | | 0.1000 | 4.0 | | 20.0 | | | |
| Benzoic acid | 0.2196 | 0.1759<br>0.2254 | 0.1584<br>0.2156 | 0.1961<br>0.2196 | 0.2111 | Ave | | 0.2027 | | | | 11.9 | | 20.0 | | | |
| Bis(2-chloroethoxy)methane | 0.4484 | 0.4618<br>0.4197 | 0.4545<br>0.3966 | 0.4730<br>0.3930 | 0.4505 | Ave | | 0.4372 | | | 0.3000 | 6.9 | | 20.0 | | | |
| 2,4-Dichlorophenol | 0.2630 | 0.2616<br>0.2479 | 0.2658<br>0.2421 | 0.2743<br>0.2367 | 0.2654 | Ave | | 0.2571 | | | 0.2000 | 5.1 | | 20.0 | | | |
| 1,2,4-Trichlorobenzene | 0.2655 | 0.3076<br>0.2480 | 0.2945<br>0.2414 | 0.2897<br>0.2370 | 0.2754 | Ave | | 0.2699 | | | | 9.7 | | 20.0 | | | |
| Naphthalene | 1.2080<br>1.0494 | 1.1687<br>0.9837 | 1.1345<br>0.9536 | 1.1160<br>0.9396 | 1.0786 | Ave | | 1.0702 | | | 0.7000 | 9.0 | | 20.0 | | | |
| 4-Chloroaniline | 0.4770 | 0.4997<br>0.4662 | 0.4663<br>0.4089 | 0.5045<br>0.4100 | 0.4730 | Ave | | 0.4632 | | | 0.0100 | 7.8 | | 20.0 | | | |
| 2,6-Dichlorophenol | 0.2525 | 0.2715<br>0.2399 | 0.2614<br>0.2321 | 0.2666<br>0.2303 | 0.2625 | Ave | | 0.2521 | | | | 6.4 | | 20.0 | | | |
| Hexachlorobutadiene | 0.1412 | 0.1455<br>0.1347 | 0.1479<br>0.1307 | 0.1551<br>0.1273 | 0.1534 | Ave | | 0.1420 | | | 0.0100 | 7.3 | | 20.0 | | | |
| Caprolactam | 0.1116 | 0.1179<br>0.1100 | 0.1187<br>0.1076 | 0.1155<br>0.1059 | 0.1083 | Ave | | 0.1119 | | | 0.0100 | 4.4 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48 Calibration End Date: 12/22/2015 20:02 Calibration ID: 32225

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 4-Chloro-3-methylphenol | 0.3396 | 0.3339<br>0.3307 | 0.3436<br>0.3159 | 0.3444<br>0.3198 | 0.3428 | Ave | | 0.3338 | | | 0.2000 | 3.3 | | 20.0 | | | |
| 2-Methylnaphthalene | 0.7542<br>0.6573 | 0.7135<br>0.6228 | 0.7088<br>0.5993 | 0.7137<br>0.5913 | 0.6889 | Ave | | 0.6722 | | | 0.4000 | 8.5 | | 20.0 | | | |
| 1-Methylnaphthalene | 0.7822<br>0.5792 | 0.6431<br>0.5405 | 0.6324<br>0.5307 | 0.6137<br>0.5229 | 0.6111 | Ave | | 0.6062 | | | | 13.1 | | 20.0 | | | |
| Hexachlorocyclopentadiene | 0.2667 | 0.2334<br>0.2581 | 0.2323<br>0.2611 | 0.2707<br>0.2532 | 0.2885 | Ave | | 0.2580 | | | 0.0500 | 7.3 | | 20.0 | | | |
| 1,2,4,5-Tetrachlorobenzene | 0.4209 | 0.4882<br>0.3822 | 0.4751<br>0.3741 | 0.4668<br>0.3558 | 0.4563 | Ave | | 0.4274 | | | 0.0100 | 12.0 | | 20.0 | | | |
| 2,4,6-Trichlorophenol | 0.3181 | 0.3464<br>0.2973 | 0.3377<br>0.2961 | 0.3259<br>0.2869 | 0.3324 | Ave | | 0.3176 | | | 0.2000 | 6.9 | | 20.0 | | | |
| 2,4,5-Trichlorophenol | 0.3295 | 0.3533<br>0.3137 | 0.3514<br>0.3103 | 0.3515<br>0.3019 | 0.3468 | Ave | | 0.3323 | | | 0.2000 | 6.4 | | 20.0 | | | |
| 1,1'-Biphenyl | 1.6646<br>1.4173 | 1.5606<br>1.3176 | 1.5365<br>1.2935 | 1.4989<br>1.2461 | 1.4773 | Ave | | 1.4458 | | | 0.0100 | 9.6 | | 20.0 | | | |
| 2-Chloronaphthalene | 1.2365<br>1.0599 | 1.1467<br>0.9906 | 1.1172<br>0.9855 | 1.1309<br>0.9466 | 1.1058 | Ave | | 1.0800 | | | 0.8000 | 8.6 | | 20.0 | | | |
| 2-Nitroaniline | 0.4791 | 0.4015<br>0.4450 | 0.4398<br>0.4446 | 0.4495<br>0.4294 | 0.4556 | Ave | | 0.4431 | | | 0.0100 | 5.0 | | 20.0 | | | |
| Dimethyl phthalate | 1.1685 | 1.2301<br>1.0811 | 1.2415<br>1.0532 | 1.2403<br>1.0143 | 1.2446 | Ave | | 1.1592 | | | 0.0100 | 8.3 | | 20.0 | | | |
| 1,3-Dinitrobenzene | 0.2038 | 0.1510<br>0.1906 | 0.1727<br>0.1928 | 0.1836<br>0.1830 | 0.2004 | Ave | | 0.1847 | | | | 9.1 | | 20.0 | | | |
| 2,6-Dinitrotoluene | 0.2845 | 0.2339<br>0.2635 | 0.2737<br>0.2627 | 0.2721<br>0.2465 | 0.2910 | Ave | | 0.2660 | | | | 7.1 | | 20.0 | | | |
| Acenaphthylene | 2.0111<br>1.8209 | 1.8589<br>1.7042 | 1.8907<br>1.6790 | 1.9261<br>1.6253 | 1.8953 | Ave | | 1.8235 | | | 0.9000 | 7.0 | | 20.0 | | | |
| 3-Nitroaniline | 0.3059 | 0.3135<br>0.3009 | 0.3010<br>0.3059 | 0.3333<br>0.2950 | 0.3242 | Ave | | 0.3099 | | | 0.0100 | 4.2 | | 20.0 | | | |
| 2,4-Dinitrophenol | 0.1315 | 0.0522<br>0.1356 | 0.0681<br>0.1409 | 0.1001<br>0.1448 | 0.1206 | Lin1 | -0.124 | 0.1424 | | | 0.0100 | | | | 0.9980 | | 0.9900 |
| 4-Nitrophenol | 0.2251 | 0.1654<br>0.2137 | 0.2048<br>0.2094 | 0.2040<br>0.2018 | 0.2088 | Ave | | 0.2041 | | | | 8.5 | | 20.0 | | | |
| Acenaphthene | 1.4039<br>1.0291 | 1.1743<br>0.9329 | 1.1711<br>0.9274 | 1.1752<br>+++++ | 1.1052 | Ave | | 1.1149 | | | 0.9000 | 14.0 | | 20.0 | | | |
| 2,4-Dinitrotoluene | 0.3447 | 0.2973<br>0.3384 | 0.3208<br>0.3455 | 0.3526<br>0.3443 | 0.3600 | Ave | | 0.3379 | | | 0.2000 | 5.9 | | 20.0 | | | |
| Dibenzofuran | 1.7448<br>1.5139 | 1.6048<br>1.4046 | 1.5927<br>1.4016 | 1.6433<br>1.3240 | 1.5827 | Ave | | 1.5347 | | | 0.8000 | 8.8 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48 Calibration End Date: 12/22/2015 20:02 Calibration ID: 32225

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2,3,4,6-Tetrachlorophenol | 0.2458 | 0.2422<br>0.2309 | 0.2652<br>0.2326 | 0.2909<br>0.2270 | 0.2863 | Ave | | 0.2526 | | | 0.0100 | 10.0 | | 20.0 | | | |
| Hexadecane | 0.9403 | 1.0869<br>0.8618 | 1.0267<br>0.8375 | 1.0377<br>0.7886 | 0.9929 | Ave | | 0.9465 | | | | 11.3 | | 20.0 | | | |
| Diethyl phthalate | 1.0291 | 1.3269<br>0.9840 | 1.2661<br>0.9888 | 1.3033<br>0.9732 | 1.1258 | Ave | | 1.1247 | | | 0.0100 | 13.6 | | 20.0 | | | |
| 4-Chlorophenyl phenyl ether | 0.5309 | 0.5565<br>0.4815 | 0.5410<br>0.4725 | 0.5691<br>0.4497 | 0.5529 | Ave | | 0.5193 | | | 0.4000 | 8.6 | | 20.0 | | | |
| 4-Nitroaniline | 0.3324 | 0.2422<br>0.3165 | 0.2947<br>0.3054 | 0.2929<br>0.2901 | 0.2629 | Ave | | 0.2921 | | | 0.0100 | 9.8 | | 20.0 | | | |
| Fluorene | 1.6545<br>1.1741 | 1.3479<br>1.0683 | 1.3255<br>+++++ | 1.3445<br>+++++ | 1.2588 | Ave | | 1.3105 | | | 0.9000 | 14.0 | | 20.0 | | | |
| 4,6-Dinitro-2-methylphenol | 0.1112 | 0.0526<br>0.1104 | 0.0698<br>0.1072 | 0.0848<br>0.1049 | 0.1057 | Lin1 | -0.065 | 0.1094 | | | 0.0100 | | | | 0.9990 | | 0.9900 |
| N-Nitrosodiphenylamine | 0.5151 | 0.5574<br>0.4402 | 0.5687<br>0.4191 | 0.5644<br>+++++ | 0.6143 | Ave | | 0.5256 | | | 0.0100 | 13.7 | | 20.0 | | | |
| 1,2-Diphenylhydrazine(as Azobenzene) | 1.1839 | 1.0932<br>1.0326 | 1.0580<br>0.9630 | 1.0800<br>0.9005 | 1.2159 | Ave | | 1.0659 | | | | 9.8 | | 20.0 | | | |
| Azobenzene | 1.1836 | 1.0962<br>1.0328 | 1.0553<br>0.9630 | 1.0792<br>0.9008 | 1.2160 | Ave | | 1.0659 | | | | 9.8 | | 20.0 | | | |
| 4-Bromophenyl phenyl ether | 0.1933 | 0.1810<br>0.1631 | 0.1883<br>0.1535 | 0.1918<br>0.1444 | 0.2103 | Ave | | 0.1782 | | | 0.1000 | 12.6 | | 20.0 | | | |
| Atrazine | 0.1767 | 0.1929<br>0.1404 | 0.1658<br>0.1377 | 0.1649<br>+++++ | 0.1931 | Ave | | 0.1673 | | | 0.0100 | 13.4 | | 20.0 | | | |
| Hexachlorobenzene | 0.2215<br>0.1812 | 0.1961<br>0.1487 | 0.1986<br>+++++ | 0.1843<br>+++++ | 0.2015 | Ave | | 0.1903 | | | 0.1000 | 11.9 | | 20.0 | | | |
| n-Octadecane | 0.6658 | 0.6723<br>0.5318 | 0.6658<br>0.4861 | 0.6500<br>+++++ | 0.7297 | Ave | | 0.6288 | | | | 13.8 | | 20.0 | | | |
| Pentachlorophenol | 0.1271 | 0.0902<br>0.1184 | 0.1054<br>0.1155 | 0.1153<br>0.1090 | 0.1297 | Ave | | 0.1138 | | | 0.0500 | 11.0 | | 20.0 | | | |
| Phenanthrene | 1.3636<br>1.0293 | 1.1644<br>0.9555 | 1.1598<br>0.9281 | 1.1515<br>0.8993 | 1.0899 | Ave | | 1.0824 | | | 0.7000 | 13.6 | | 20.0 | | | |
| Anthracene | 1.2554<br>1.0489 | 1.1608<br>0.9736 | 1.1511<br>0.9503 | 1.1500<br>0.9134 | 1.1164 | Ave | | 1.0800 | | | 0.7000 | 10.6 | | 20.0 | | | |
| Carbazole | 1.0373 | 1.0814<br>0.8833 | 1.0489<br>0.8362 | 0.8963<br>0.7816 | 0.9925 | Ave | | 0.9447 | | | 0.0100 | 11.7 | | 20.0 | | | |
| Di-n-butyl phthalate | 1.4091 | 1.1737<br>1.1938 | 0.9563<br>1.1660 | 1.0751<br>1.1386 | 1.4093 | Ave | | 1.1902 | | | 0.0100 | 13.0 | | 20.0 | | | |
| Fluoranthene | 1.2783<br>1.1573 | 1.1582<br>1.0065 | 1.0063<br>0.9626 | 0.9440<br>0.9144 | 1.1785 | Ave | | 1.0674 | | | 0.6000 | 12.0 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48 Calibration End Date: 12/22/2015 20:02 Calibration ID: 32225

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Benzidine | 0.5165 | +++++<br>0.5681 | 0.2518<br>0.5466 | 0.2776<br>0.5590 | 0.4607 | Lin1 | -0.847 | 0.5726 | | | | | | | 0.9980 | | 0.9900 |
| Pyrene | 1.3761<br>1.0608 | 1.1093<br>0.9884 | 1.0001<br>0.9683 | 1.0079<br>0.9541 | 1.0427 | Ave | | 1.0564 | | | 0.6000 | 12.2 | | 20.0 | | | |
| Butyl benzyl phthalate | 0.6006 | 0.5689<br>0.5758 | 0.5694<br>0.5622 | 0.5753<br>0.5603 | 0.5729 | Ave | | 0.5732 | | | 0.0100 | 2.2 | | 20.0 | | | |
| Bis(2-ethylhexyl) phthalate | 0.8276 | 0.7165<br>0.7886 | 0.7847<br>0.7798 | 0.7864<br>0.7843 | 0.8004 | Ave | | 0.7835 | | | 0.0100 | 4.0 | | 20.0 | | | |
| 3,3'-Dichlorobenzidine | 0.2713 | 0.2979<br>0.2708 | 0.3205<br>0.2807 | 0.3240<br>0.2773 | 0.2529 | Ave | | 0.2869 | | | 0.0100 | 8.8 | | 20.0 | | | |
| Benzo[a]anthracene | 1.2777<br>1.0264 | 1.0668<br>0.9842 | 1.0996<br>0.9616 | 1.0823<br>0.9473 | 1.0505 | Ave | | 1.0552 | | | 0.8000 | 9.4 | | 20.0 | | | |
| Chrysene | 1.2706<br>0.9658 | 1.0700<br>0.9125 | 1.0175<br>0.8900 | 1.0526<br>0.8863 | 0.9943 | Ave | | 1.0066 | | | 0.7000 | 11.9 | | 20.0 | | | |
| Di-n-octyl phthalate | 1.6857 | 1.4216<br>1.6330 | 1.6580<br>1.5940 | 1.7040<br>1.5678 | 1.6935 | Ave | | 1.6197 | | | 0.0100 | 5.8 | | 20.0 | | | |
| Benzo[b]fluoranthene | 1.1807<br>1.1236 | 1.0820<br>1.0918 | 1.1664<br>1.0534 | 1.1835<br>1.0502 | 1.1488 | Ave | | 1.1200 | | | 0.7000 | 4.7 | | 20.0 | | | |
| Benzo[k]fluoranthene | 1.4389<br>1.1869 | 1.2148<br>1.0837 | 1.2153<br>1.0932 | 1.2694<br>1.0315 | 1.1923 | Ave | | 1.1918 | | | 0.7000 | 10.1 | | 20.0 | | | |
| Benzo[a]pyrene | 1.1499<br>1.1196 | 1.0799<br>1.0664 | 1.1335<br>1.0496 | 1.1666<br>1.0260 | 1.1404 | Ave | | 1.1035 | | | 0.7000 | 4.5 | | 20.0 | | | |
| Indeno[1,2,3-cd]pyrene | 1.1649<br>1.1153 | 1.0698<br>1.0519 | 1.1429<br>1.0363 | 1.1927<br>1.0101 | 1.1379 | Ave | | 1.1024 | | | 0.5000 | 5.7 | | 20.0 | | | |
| Dibenz(a,h)anthracene | 0.7391<br>0.9367 | 0.9194<br>0.8858 | 0.9181<br>0.8521 | 0.9713<br>0.8483 | 0.9473 | Ave | | 0.8909 | | | 0.4000 | 7.9 | | 20.0 | | | |
| Benzo[g,h,i]perylene | 1.1015<br>0.9455 | 0.9093<br>0.9004 | 0.9938<br>0.8941 | 1.0092<br>0.8761 | 0.9705 | Ave | | 0.9556 | | | 0.5000 | 7.5 | | 20.0 | | | |
| 2-Fluorophenol (Surr) | 1.4614 | 1.3531<br>1.3811 | 1.4159<br>1.3960 | 1.4306<br>1.3999 | 1.3875 | Ave | | 1.4032 | | | | 2.3 | | 20.0 | | | |
| Phenol-d5 (Surr) | 1.8754 | 1.8220<br>1.7435 | 1.9269<br>1.7574 | 1.8860<br>1.7393 | 1.8389 | Ave | | 1.8237 | | | | 3.9 | | 20.0 | | | |
| Nitrobenzene-d5 (Surr) | 0.5258<br>0.4815 | 0.4786<br>0.4664 | 0.5039<br>0.4575 | 0.5053<br>0.4492 | 0.4832 | Ave | | 0.4835 | | | | 5.1 | | 20.0 | | | |
| 2-Fluorobiphenyl (Surr) | 1.4586<br>1.1666 | 1.3120<br>1.0886 | 1.2881<br>1.0717 | 1.2857<br>1.0380 | 1.2313 | Ave | | 1.2156 | | | | 11.2 | | 20.0 | | | |
| 2,4,6-Tribromophenol (Surr) | 0.1215 | 0.1020<br>0.1155 | 0.1273<br>0.1171 | 0.1361<br>0.1179 | 0.1317 | Ave | | 0.1211 | | | | 8.8 | | 20.0 | | | |
| Terphenyl-d14 (Surr) | 0.9148<br>0.7155 | 0.7332<br>0.6735 | 0.7562<br>0.6547 | 0.7212<br>0.6546 | 0.7286 | Ave | | 0.7280 | | | | 10.8 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48 Calibration End Date: 12/22/2015 20:02 Calibration ID: 32225

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-211933/6 | 51222006.D |
| Level 2 | STD2 240-211933/5 | 51222005.D |
| Level 3 | STD3 240-211933/4 | 51222004.D |
| Level 4 | STD4 240-211933/3 | 51222003.D |
| Level 5 | STD5 240-211933/2 | 51222002.D |
| Level 6 | STD6 240-211933/10 | 51222010.D |
| Level 7 | STD7 240-211933/9 | 51222009.D |
| Level 8 | STD8 240-211933/8 | 51222008.D |
| Level 9 | STD9 240-211933/7 | 51222007.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1,4-Dioxane | DCB | Ave | 449195 | 22401<br>659966 | 45106<br>875199 | 86174<br>1009112 | 210938 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| N-Nitrosodimethylamine | DCB | Ave | 680038 | 32992<br>998454 | 62682<br>1297577 | 130777<br>1514764 | 301281 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Pyridine | DCB | Ave | 1202602 | 59903<br>1788998 | 120933<br>2282211 | 224253<br>2660177 | 534780 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzaldehyde | DCB | Ave | 1543587 | 98802<br>2484066 | 187185<br>2991393 | 387036<br>3335858 | 851877 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Phenol | DCB | Ave | 1214360 | 68676<br>1865920 | 135337<br>2303476 | 266289<br>2639426 | 608595 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Aniline | DCB | Ave | 1532152 | 84892<br>2301384 | 159599<br>2883535 | 324167<br>3348828 | 748775 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Bis(2-chloroethyl)ether | DCB | Ave | 1035534 | 54111<br>1559640 | 106100<br>1931236 | 219807<br>2245739 | 516571 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Chlorophenol | DCB | Ave | 870389 | 46290<br>1295178 | 93532<br>1621907 | 185526<br>1857905 | 436804 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| n-Decane | DCB | Ave | 1129120 | 60007<br>1642170 | 120508<br>2038274 | 245431<br>2320206 | 549881 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,3-Dichlorobenzene | DCB | Ave | 915628 | 52705<br>1368158 | 102451<br>1711292 | 197194<br>1948583 | 463873 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,4-Dichlorobenzene | DCB | Ave | 906737 | 50244<br>1382571 | 99724<br>1719256 | 201698<br>1978465 | 459953 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzyl alcohol | DCB | Ave | 654233 | 37559<br>1002672 | 68629<br>1227342 | 140774<br>1442611 | 325311 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2-Dichlorobenzene | DCB | Ave | 880477 | 50578<br>1308977 | 96124<br>1641132 | 195718<br>1852472 | 447437 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Methylphenol | DCB | Ave | 896896 | 48466<br>1361792 | 97165<br>1694511 | 193788<br>1944423 | 445264 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48 Calibration End Date: 12/22/2015 20:02 Calibration ID: 32225

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| bis (2-chloroisopropyl) ether | DCB | Ave | 1285204 | 71162<br>1936692 | 136330<br>2385690 | 270279<br>2737528 | 630107 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Indene | DCB | Ave | 3030051 | 181779<br>4454751 | 349215<br>5475325 | 679786<br>6235361 | 1563344 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 3 & 4 Methylphenol | DCB | Ave | 944263 | 49082<br>1428535 | 100430<br>1756740 | 201580<br>2024340 | 474193 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| N-Nitrosodi-n-propylamine | DCB | Ave | 731056 | 42299<br>1114709 | 74891<br>1354073 | 159494<br>1602839 | 377394 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Acetophenone | DCB | Ave | 1309809 | 76240<br>2011573 | 140926<br>2441181 | 287028<br>2827168 | 672275 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachloroethane | DCB | Ave | 378039 | 22014<br>575981 | 42268<br>723880 | 82971<br>824648 | 192399 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Nitrobenzene | NPT | Ave | 1147098 | 60833<br>1744452 | 117024<br>2147390 | 244647<br>2517728 | 567391 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Isophorone | NPT | Ave | 2036234 | 108717<br>3095872 | 209731<br>3766245 | 437472<br>4417157 | 1024432 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dimethylphenol | NPT | Ave | 936317 | 51825<br>1431456 | 102596<br>1722767 | 206887<br>1993284 | 489018 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Nitrophenol | NPT | Ave | 410869 | 20266<br>618816 | 40127<br>744078 | 84813<br>891093 | 205322 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzoic acid | NPT | Ave | 1070859 | 45015<br>1754659 | 78893<br>2110188 | 197692<br>2539902 | 516472 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Bis(2-chloroethoxy)methane | NPT | Ave | 1093511 | 59103<br>1633434 | 113224<br>1940878 | 238459<br>2272737 | 551115 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dichlorophenol | NPT | Ave | 641385 | 33475<br>964596 | 66205<br>1184923 | 138292<br>1368931 | 324742 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2,4-Trichlorobenzene | NPT | Ave | 647365 | 39366<br>965232 | 73354<br>1181347 | 146020<br>1370743 | 336916 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Naphthalene | NPT | Ave | 31013<br>2559045 | 149574<br>3828054 | 282596<br>4667248 | 562552<br>5434110 | 1319584 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Chloroaniline | NPT | Ave | 1163154 | 63950<br>1814156 | 116152<br>2001389 | 254303<br>2370956 | 578603 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,6-Dichlorophenol | NPT | Ave | 615862 | 34746<br>933552 | 65122<br>1136080 | 134374<br>1331798 | 321081 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorobutadiene | NPT | Ave | 344216 | 18627<br>524116 | 36852<br>639818 | 78168<br>735954 | 187656 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Caprolactam | NPT | Ave | 544125 | 30186<br>856410 | 59153<br>1053173 | 116475<br>1225368 | 264882 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 4-Chloro-3-methylphenol | NPT | Ave | 828150 | 42735<br>1286780 | 85600<br>1546019 | 173589<br>1849173 | 419398 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Methylnaphthalene | NPT | Ave | 19362<br>1602796 | 91309<br>2423525 | 176560<br>2933099 | 359794<br>3419662 | 842786 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48 Calibration End Date: 12/22/2015 20:02 Calibration ID: 32225

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 1-Methylnaphthalene | NPT | Ave | 20082<br>1412414 | 82310<br>2103465 | 157535<br>2597181 | 309390<br>3024116 | 747599 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexachlorocyclopentadiene | ANT | Ave | <br>341214 | 15917<br>529599 | 30919<br>655199 | 72390<br>772405 | 185901 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,2,4,5-Tetrachlorobenzene | ANT | Ave | <br>538528 | 33294<br>784302 | 63236<br>938932 | 124816<br>1085449 | 294024 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,6-Trichlorophenol | ANT | Ave | <br>407075 | 23626<br>610049 | 44951<br>743055 | 87130<br>875250 | 214207 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,5-Trichlorophenol | ANT | Ave | <br>421575 | 24099<br>643684 | 46771<br>778743 | 93982<br>920956 | 223457 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,1'-Biphenyl | ANT | Ave | 22687<br>1813500 | 106434<br>2704010 | 204525<br>3246303 | 400765<br>3801339 | 951866 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Chloronaphthalene | ANT | Ave | 16852<br>1356192 | 78211<br>2033040 | 148715<br>2473224 | 302381<br>2887651 | 712497 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Nitroaniline | ANT | Ave | <br>613070 | 27382<br>913144 | 58544<br>1115704 | 120196<br>1309891 | 293570 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dimethyl phthalate | ANT | Ave | <br>1495160 | 83898<br>2218721 | 165259<br>2643222 | 331618<br>3094270 | 801971 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 1,3-Dinitrobenzene | ANT | Ave | <br>260758 | 10299<br>391173 | 22990<br>483898 | 49095<br>558377 | 129126 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,6-Dinitrotoluene | ANT | Ave | <br>364023 | 15953<br>540839 | 36427<br>659392 | 72763<br>751876 | 187499 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Acenaphthylene | ANT | Ave | 27410<br>2329849 | 126782<br>3497456 | 251680<br>4213699 | 514984<br>4958173 | 1221242 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 3-Nitroaniline | ANT | Ave | <br>391461 | 21379<br>617518 | 40062<br>767616 | 89108<br>899942 | 208872 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4-Dinitrophenol | ANT | Lin1 | <br>336456 | 7119<br>556645 | 18124<br>707053 | 53531<br>883368 | 155411 | <br>20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| 4-Nitrophenol | ANT | Ave | <br>576019 | 22560<br>877039 | 54511<br>1051021 | 109100<br>1231049 | 269023 | <br>20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Acenaphthene | ANT | Ave | 19134<br>1316697 | 80092<br>1914552 | 155888<br>2327390 | 314209<br>+++++ | 712118 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>+++++ | 5.00 |
| 2,4-Dinitrotoluene | ANT | Ave | <br>440986 | 20280<br>694467 | 42699<br>867058 | 94277<br>1050304 | 231966 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dibenzofuran | ANT | Ave | 23780<br>1937015 | 109450<br>2882574 | 212014<br>3517678 | 439371<br>4038942 | 1019817 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,3,4,6-Tetrachlorophenol | ANT | Ave | <br>314505 | 16516<br>473830 | 35304<br>583802 | 77780<br>692340 | 184485 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Hexadecane | ANT | Ave | <br>1203112 | 74130<br>1768569 | 136663<br>2101818 | 277447<br>2405620 | 639759 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Diethyl phthalate | ANT | Ave | <br>1316809 | 90501<br>2019395 | 168538<br>2481446 | 348481<br>2968937 | 725401 | <br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48 Calibration End Date: 12/22/2015 20:02 Calibration ID: 32225

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 4-Chlorophenyl phenyl ether | ANT | Ave | 679329 | 37955<br>988084 | 72010<br>1185912 | 152162<br>1371994 | 356247 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Nitroaniline | ANT | Ave | 425352 | 16518<br>649627 | 39222<br>766343 | 78313<br>884950 | 169390 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Fluorene | ANT | Ave | 22550<br>1502232 | 91934<br>2192504 | 176439<br>+++++ | 359497<br>+++++ | 811097 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>+++++ | 2.00<br>+++++ | 5.00 |
| 4,6-Dinitro-2-methylphenol | PHN | Lin1 | 415420 | 11580<br>691700 | 30020<br>850540 | 74400<br>1031899 | 196586 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| N-Nitrosodiphenylamine | PHN | Ave | 1924208 | 122742<br>2758099 | 244597<br>3324572 | 495375<br>+++++ | 1142182 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>+++++ | 10.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | PHN | Ave | 2211164 | 120365<br>3234768 | 227501<br>3819938 | 473981<br>4428921 | 1130415 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Azobenzene | PHN | Ave | 2210700 | 120695<br>3235201 | 226933<br>3819938 | 473601<br>4430238 | 1130519 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 4-Bromophenyl phenyl ether | PHN | Ave | 361033 | 19935<br>510826 | 40493<br>608824 | 84168<br>710178 | 195539 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Atrazine | PHN | Ave | 659970 | 42469<br>879531 | 71290<br>1092522 | 144715<br>+++++ | 359116 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>+++++ | 10.0 |
| Hexachlorobenzene | PHN | Ave | 4862<br>338463 | 21595<br>465671 | 42714<br>+++++ | 80869<br>+++++ | 187374 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>+++++ | 2.00<br>+++++ | 5.00 |
| n-Octadecane | PHN | Ave | 1243629 | 74020<br>1665745 | 143164<br>1928391 | 285258<br>+++++ | 678442 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>+++++ | 5.00 |
| Pentachlorophenol | PHN | Ave | 474667 | 19864<br>742055 | 45340<br>916603 | 101212<br>1072301 | 241174 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Phenanthrene | PHN | Ave | 29927<br>1922443 | 128213<br>2993255 | 249404<br>3681414 | 505368<br>4422963 | 1013305 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Anthracene | PHN | Ave | 27551<br>1959122 | 127809<br>3049751 | 247523<br>3769753 | 504697<br>4491990 | 1037906 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Carbazole | PHN | Ave | 1937441 | 119071<br>2766965 | 225547<br>3317035 | 393358<br>3844090 | 922727 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Di-n-butyl phthalate | PHN | Ave | 2631873 | 129238<br>3739726 | 205630<br>4625358 | 471802<br>5599721 | 1310292 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Fluoranthene | PHN | Ave | 28055<br>2161544 | 127532<br>3152930 | 216389<br>3818462 | 414273<br>4497246 | 1095671 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzidine | CRY | Lin1 | 2187463 | +++++<br>3793518 | 109369<br>4524072 | 247197<br>5523609 | 1034312 | 20.0 | +++++<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Pyrene | CRY | Ave | 30318<br>2246219 | 123771<br>3299988 | 217225<br>4007174 | 448691<br>4714311 | 1170469 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Butyl benzyl phthalate | CRY | Ave | 1271886 | 63471<br>1922532 | 123670<br>2326509 | 256105<br>2768388 | 643093 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Bis(2-ethylhexyl) phthalate | CRY | Ave | 1752462 | 79944<br>2632974 | 170449<br>3226911 | 350104<br>3875236 | 898434 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 211933

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/22/2015 12:48 Calibration End Date: 12/22/2015 20:02 Calibration ID: 32225

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 |
| 3,3'-Dichlorobenzidine | CRY | Ave | 1149020 | 66487<br>1808538 | 139250<br>2323310 | 288497<br>2740041 | 567781 | 20.0 | 1.00<br>30.0 | 2.00<br>40.0 | 4.00<br>50.0 | 10.0 |
| Benzo[a]anthracene | CRY | Ave | 28151<br>2173411 | 119029<br>3286100 | 238855<br>3979280 | 481831<br>4680672 | 1179191 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Chrysene | CRY | Ave | 27993<br>2045001 | 119388<br>3046622 | 221013<br>3683059 | 468568<br>4379009 | 1116152 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Di-n-octyl phthalate | PRY | Ave | 3071125 | 132042<br>4783871 | 292269<br>5849192 | 618688<br>6969419 | 1574787 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[b]fluoranthene | PRY | Ave | 22149<br>2046981 | 100497<br>3198445 | 205606<br>3865380 | 429694<br>4668631 | 1068304 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[k]fluoranthene | PRY | Ave | 26992<br>2162415 | 112836<br>3174639 | 214232<br>4011475 | 460873<br>4585294 | 1108725 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[a]pyrene | PRY | Ave | 21570<br>2039735 | 100302<br>3123991 | 199820<br>3851450 | 423553<br>4560949 | 1060514 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Indeno[1,2,3-cd]pyrene | PRY | Ave | 21851<br>2031848 | 99364<br>3081525 | 201467<br>3802782 | 433046<br>4490199 | 1058176 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Dibenz(a,h)anthracene | PRY | Ave | 13864<br>1706509 | 85396<br>2594848 | 161846<br>3126710 | 352670<br>3770966 | 880949 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Benzo[g,h,i]perylene | PRY | Ave | 20662<br>1722480 | 84458<br>2637507 | 175189<br>3280702 | 366415<br>3894742 | 902525 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Fluorophenol (Surr) | DCB | Ave | 876049 | 44324<br>1338707 | 88355<br>1673378 | 180126<br>1939453 | 420218 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Phenol-d5 (Surr) | DCB | Ave | 1124238 | 59683<br>1689949 | 120239<br>2106588 | 237475<br>2409627 | 556917 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Nitrobenzene-d5 (Surr) | NPT | Ave | 13498<br>1174187 | 61250<br>1815222 | 125526<br>2238915 | 254715<br>2597908 | 591164 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2-Fluorobiphenyl (Surr) | ANT | Ave | 19879<br>1492624 | 89485<br>2234027 | 171462<br>2689502 | 343758<br>3166626 | 793353 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| 2,4,6-Tribromophenol (Surr) | ANT | Ave | 155491 | 6960<br>237109 | 16939<br>293777 | 36403<br>359708 | 84858 | 10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |
| Terphenyl-d14 (Surr) | CRY | Ave | 20154<br>1515158 | 81806<br>2248658 | 164256<br>2709388 | 321041<br>3234339 | 817844 | 0.100<br>10.0 | 0.500<br>15.0 | 1.00<br>20.0 | 2.00<br>25.0 | 5.00 |

Curve Type Legend:

| |
|---|
| Ave = Average ISTD |
| Lin1 = Linear 1/conc ISTD |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID (1): CCV 240-212723/2 CCVIS Instrument ID (1): A4HP9

GC Column (1): RXI-5SILMS ID: 0.45(mm) Date Analyzed (1): 12/31/2015 09:36

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 2-Fluorophenol (Surr) | 5.24 | 100.0 |
| Benzaldehyde | 6.07 | 100.0 |
| Phenol-d5 (Surr) | 6.10 | 100.0 |
| Phenol | 6.11 | 100.0 |
| Bis(2-chloroethyl)ether | 6.20 | 100.0 |
| 2-Chlorophenol | 6.28 | 100.0 |
| 2-Methylphenol | 6.65 | 100.0 |
| bis (2-chloroisopropyl) ether | 6.68 | 100.0 |
| 3 & 4 Methylphenol | 6.78 | 100.0 |
| N-Nitrosodi-n-propylamine | 6.79 | 100.0 |
| Acetophenone | 6.81 | 100.0 |
| Hexachloroethane | 6.93 | 100.0 |
| Nitrobenzene-d5 (Surr) | 6.95 | 100.0 |
| Nitrobenzene | 6.96 | 100.0 |
| Isophorone | 7.16 | 100.0 |
| 2-Nitrophenol | 7.24 | 100.0 |
| 2,4-Dimethylphenol | 7.24 | 100.0 |
| Bis(2-chloroethoxy)methane | 7.32 | 100.0 |
| 2,4-Dichlorophenol | 7.45 | 100.0 |
| Naphthalene | 7.60 | 100.0 |
| 4-Chloroaniline | 7.62 | 100.0 |
| Hexachlorobutadiene | 7.70 | 100.0 |
| Caprolactam | 7.91 | 100.0 |
| 4-Chloro-3-methylphenol | 8.01 | 100.0 |
| 2-Methylnaphthalene | 8.19 | 100.0 |
| Hexachlorocyclopentadiene | 8.33 | 100.0 |
| 2,4,6-Trichlorophenol | 8.42 | 100.0 |
| 2,4,5-Trichlorophenol | 8.46 | 100.0 |
| 2-Fluorobiphenyl (Surr) | 8.48 | 100.0 |
| 1,1'-Biphenyl | 8.58 | 100.0 |
| 2-Chloronaphthalene | 8.61 | 100.0 |
| 2-Nitroaniline | 8.67 | 100.0 |
| Dimethyl phthalate | 8.80 | 100.0 |
| 2,6-Dinitrotoluene | 8.86 | 100.0 |
| Acenaphthylene | 8.97 | 100.0 |
| 3-Nitroaniline | 9.01 | 100.0 |
| 2,4-Dinitrophenol | 9.10 | 100.0 |
| Acenaphthene | 9.12 | 100.0 |
| 4-Nitrophenol | 9.12 | 100.0 |
| 2,4-Dinitrotoluene | 9.21 | 100.0 |
| Dibenzofuran | 9.26 | 100.0 |
| Diethyl phthalate | 9.39 | 100.0 |
| 4-Chlorophenyl phenyl ether | 9.52 | 100.0 |
| 4-Nitroaniline | 9.54 | 100.0 |
| Fluorene | 9.55 | 100.0 |

FORM VI
RESOLUTION CHECK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID (1): CCV 240-212723/2 CCVIS Instrument ID (1): A4HP9

GC Column (1): RXI-5SILMS ID: 0.45(mm) Date Analyzed (1): 12/31/2015 09:36

| ANALYTE | RT | RESOLUTION (%) |
|---|---|---|
| 4,6-Dinitro-2-methylphenol | 9.56 | 100.0 |
| N-Nitrosodiphenylamine | 9.62 | 100.0 |
| 2,4,6-Tribromophenol (Surr) | 9.76 | 100.0 |
| 4-Bromophenyl phenyl ether | 9.94 | 100.0 |
| Hexachlorobenzene | 10.04 | 100.0 |
| Atrazine | 10.04 | 100.0 |
| Pentachlorophenol | 10.19 | 100.0 |
| Phenanthrene | 10.38 | 100.0 |
| Anthracene | 10.42 | 100.0 |
| Carbazole | 10.54 | 100.0 |
| Di-n-butyl phthalate | 10.76 | 100.0 |
| Fluoranthene | 11.40 | 100.0 |
| Pyrene | 11.60 | 100.0 |
| Terphenyl-d14 (Surr) | 11.67 | 100.0 |
| Butyl benzyl phthalate | 12.04 | 100.0 |
| Bis(2-ethylhexyl) phthalate | 12.53 | 100.0 |
| 3,3'-Dichlorobenzidine | 12.58 | 100.0 |
| Benzo[a]anthracene | 12.65 | 100.0 |
| Chrysene | 12.69 | 100.0 |
| Di-n-octyl phthalate | 13.21 | 100.0 |
| Benzo[b]fluoranthene | 13.94 | 26.50 |
| Benzo[k]fluoranthene | 13.98 | 100.0 |
| Benzo[a]pyrene | 14.44 | 100.0 |
| Indeno[1,2,3-cd]pyrene | 16.56 | 100.0 |
| Dibenz(a,h)anthracene | 16.57 | 100.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-211933/11 Calibration Date: 12/22/2015 16:38

Instrument ID: A4HP9 Calib Start Date: 12/22/2015 12:48

GC Column: RXI-5SILMS ID: 0.45(mm) Calib End Date: 12/22/2015 20:02

Lab File ID: 51222011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.7095 | 0.6907 | | 9.73 | 10.0 | -2.7 | 30.0 |
| N-Nitrosodimethylamine | Ave | 1.048 | 0.9229 | | 8.80 | 10.0 | -12.0 | 30.0 |
| Pyridine | Ave | 1.874 | 1.856 | | 9.90 | 10.0 | -1.0 | 30.0 |
| Benzaldehyde | Ave | 1.371 | 1.338 | 0.0100 | 19.5 | 20.0 | -2.4 | 30.0 |
| Phenol | Ave | 2.021 | 1.930 | 0.8000 | 9.55 | 10.0 | -4.5 | 30.0 |
| Aniline | Ave | 2.494 | 2.398 | | 9.62 | 10.0 | -3.8 | 30.0 |
| Bis(2-chloroethyl)ether | Ave | 1.672 | 1.620 | 0.7000 | 9.69 | 10.0 | -3.1 | 30.0 |
| 2-Chlorophenol | Ave | 1.414 | 1.354 | 0.8000 | 9.58 | 10.0 | -4.2 | 30.0 |
| n-Decane | Ave | 1.810 | 1.726 | | 9.53 | 10.0 | -4.7 | 30.0 |
| 1,3-Dichlorobenzene | Ave | 1.515 | 1.414 | | 9.33 | 10.0 | -6.7 | 30.0 |
| 1,4-Dichlorobenzene | Ave | 1.507 | 1.438 | | 9.54 | 10.0 | -4.6 | 30.0 |
| Benzyl alcohol | Ave | 1.079 | 1.039 | | 9.64 | 10.0 | -3.6 | 30.0 |
| 1,2-Dichlorobenzene | Ave | 1.455 | 1.368 | | 9.40 | 10.0 | -6.0 | 30.0 |
| 2-Methylphenol | Ave | 1.471 | 1.450 | 0.7000 | 9.86 | 10.0 | -1.4 | 30.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.087 | 2.036 | | 9.76 | 10.0 | -2.4 | 30.0 |
| Indene | Ave | 2.527 | 2.240 | | 17.7 | 20.0 | -11.3 | 30.0 |
| 3 & 4 Methylphenol | Ave | 1.531 | 1.501 | | 9.80 | 10.0 | -2.0 | 30.0 |
| N-Nitrosodi-n-propylamine | Ave | 1.208 | 1.181 | 0.5000 | 9.78 | 10.0 | -2.2 | 30.0 |
| Acetophenone | Ave | 2.178 | 2.099 | 0.0100 | 9.64 | 10.0 | -3.6 | 30.0 |
| Hexachloroethane | Ave | 0.6335 | 0.6019 | 0.3000 | 9.50 | 10.0 | -5.0 | 30.0 |
| Nitrobenzene | Ave | 0.4609 | 0.4393 | 0.2000 | 9.53 | 10.0 | -4.7 | 30.0 |
| Isophorone | Ave | 0.8201 | 0.8061 | 0.4000 | 9.83 | 10.0 | -1.7 | 30.0 |
| 2,4-Dimethylphenol | Ave | 0.3844 | 0.3558 | 0.2000 | 9.25 | 10.0 | -7.5 | 30.0 |
| 2-Nitrophenol | Ave | 0.1611 | 0.1577 | 0.1000 | 9.78 | 10.0 | -2.2 | 30.0 |
| Benzoic acid | Ave | 0.2027 | 0.2184 | | 21.6 | 20.0 | 7.8 | 30.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4372 | 0.4329 | 0.3000 | 9.90 | 10.0 | -1.0 | 30.0 |
| 2,4-Dichlorophenol | Ave | 0.2571 | 0.2384 | 0.2000 | 9.27 | 10.0 | -7.3 | 30.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.2699 | 0.2492 | | 9.23 | 10.0 | -7.7 | 30.0 |
| Naphthalene | Ave | 1.070 | 0.9806 | 0.7000 | 9.16 | 10.0 | -8.4 | 30.0 |
| 4-Chloroaniline | Ave | 0.4632 | 0.4311 | 0.0100 | 9.31 | 10.0 | -6.9 | 30.0 |
| 2,6-Dichlorophenol | Ave | 0.2521 | 0.2384 | | 9.46 | 10.0 | -5.4 | 30.0 |
| Hexachlorobutadiene | Ave | 0.1420 | 0.1370 | 0.0100 | 9.65 | 10.0 | -3.5 | 30.0 |
| Caprolactam | Ave | 0.1119 | 0.1064 | 0.0100 | 19.0 | 20.0 | -5.0 | 30.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3338 | 0.3264 | 0.2000 | 9.78 | 10.0 | -2.2 | 30.0 |
| 2-Methylnaphthalene | Ave | 0.6722 | 0.5948 | 0.4000 | 8.85 | 10.0 | -11.5 | 30.0 |
| 1-Methylnaphthalene | Ave | 0.6062 | 0.5689 | | 9.39 | 10.0 | -6.1 | 30.0 |
| Hexachlorocyclopentadiene | Ave | 0.2580 | 0.2608 | 0.0500 | 10.1 | 10.0 | 1.1 | 30.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4274 | 0.3959 | 0.0100 | 9.26 | 10.0 | -7.4 | 30.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3176 | 0.3094 | 0.2000 | 9.74 | 10.0 | -2.6 | 30.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3323 | 0.3342 | 0.2000 | 10.1 | 10.0 | 0.6 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-211933/11 Calibration Date: 12/22/2015 16:38

Instrument ID: A4HP9 Calib Start Date: 12/22/2015 12:48

GC Column: RXI-5SILMS ID: 0.45(mm) Calib End Date: 12/22/2015 20:02

Lab File ID: 51222011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.446 | 1.344 | 0.0100 | 9.29 | 10.0 | -7.1 | 30.0 |
| 2-Chloronaphthalene | Ave | 1.080 | 1.000 | 0.8000 | 9.26 | 10.0 | -7.4 | 30.0 |
| 2-Nitroaniline | Ave | 0.4431 | 0.4567 | 0.0100 | 10.3 | 10.0 | 3.1 | 30.0 |
| Dimethyl phthalate | Ave | 1.159 | 1.127 | 0.0100 | 9.72 | 10.0 | -2.8 | 30.0 |
| 1,3-Dinitrobenzene | Ave | 0.1847 | 0.1908 | | 10.3 | 10.0 | 3.3 | 30.0 |
| 2,6-Dinitrotoluene | Ave | 0.2660 | 0.2738 | | 10.3 | 10.0 | 2.9 | 30.0 |
| Acenaphthylene | Ave | 1.824 | 1.707 | 0.9000 | 9.36 | 10.0 | -6.4 | 30.0 |
| 3-Nitroaniline | Ave | 0.3099 | 0.3008 | 0.0100 | 9.70 | 10.0 | -3.0 | 30.0 |
| 2,4-Dinitrophenol | Lin1 | | 0.1273 | 0.0100 | 18.8 | 20.0 | -6.2 | 30.0 |
| 4-Nitrophenol | Ave | 0.2041 | 0.2106 | | 20.6 | 20.0 | 3.2 | 30.0 |
| Acenaphthene | Ave | 1.115 | 0.9914 | 0.9000 | 8.89 | 10.0 | -11.1 | 30.0 |
| 2,4-Dinitrotoluene | Ave | 0.3379 | 0.3351 | 0.2000 | 9.91 | 10.0 | -0.9 | 30.0 |
| Dibenzofuran | Ave | 1.535 | 1.447 | 0.8000 | 9.43 | 10.0 | -5.7 | 30.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.2526 | 0.2366 | 0.0100 | 9.37 | 10.0 | -6.3 | 30.0 |
| Hexadecane | Ave | 0.9465 | 0.9020 | | 9.53 | 10.0 | -4.7 | 30.0 |
| Diethyl phthalate | Ave | 1.125 | 1.003 | 0.0100 | 8.92 | 10.0 | -10.8 | 30.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.5193 | 0.4974 | 0.4000 | 9.58 | 10.0 | -4.2 | 30.0 |
| 4-Nitroaniline | Ave | 0.2921 | 0.3124 | 0.0100 | 10.7 | 10.0 | 6.9 | 30.0 |
| Fluorene | Ave | 1.311 | 1.120 | 0.9000 | 8.55 | 10.0 | -14.5 | 30.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1104 | 0.0100 | 20.8 | 20.0 | 3.8 | 30.0 |
| N-Nitrosodiphenylamine | Ave | 0.5256 | 0.4764 | 0.0100 | 18.1 | 20.0 | -9.4 | 30.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 1.066 | 1.169 | | 11.0 | 10.0 | 9.7 | 30.0 |
| Azobenzene | Ave | 1.066 | 1.169 | | 11.0 | 10.0 | 9.7 | 30.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.1782 | 0.1789 | 0.1000 | 10.0 | 10.0 | 0.4 | 30.0 |
| Atrazine | Ave | 0.1673 | 0.1589 | 0.0100 | 19.0 | 20.0 | -5.1 | 30.0 |
| Hexachlorobenzene | Ave | 0.1903 | 0.1686 | 0.1000 | 8.86 | 10.0 | -11.4 | 30.0 |
| n-Octadecane | Ave | 0.6288 | 0.6223 | | 9.90 | 10.0 | -1.0 | 30.0 |
| Pentachlorophenol | Ave | 0.1138 | 0.1256 | 0.0500 | 22.1 | 20.0 | 10.4 | 30.0 |
| Phenanthrene | Ave | 1.082 | 0.9592 | 0.7000 | 8.86 | 10.0 | -11.4 | 30.0 |
| Anthracene | Ave | 1.080 | 0.9938 | 0.7000 | 9.20 | 10.0 | -8.0 | 30.0 |
| Carbazole | Ave | 0.9447 | 0.9472 | 0.0100 | 10.0 | 10.0 | 0.3 | 30.0 |
| Di-n-butyl phthalate | Ave | 1.190 | 1.342 | 0.0100 | 11.3 | 10.0 | 12.8 | 30.0 |
| Fluoranthene | Ave | 1.067 | 1.092 | 0.6000 | 10.2 | 10.0 | 2.3 | 30.0 |
| Benzidine | Lin1 | | 0.5460 | | 20.5 | 20.0 | 2.7 | 30.0 |
| Pyrene | Ave | 1.056 | 0.999 | 0.6000 | 9.46 | 10.0 | -5.4 | 30.0 |
| Butyl benzyl phthalate | Ave | 0.5732 | 0.5786 | 0.0100 | 10.1 | 10.0 | 0.9 | 30.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.7835 | 0.7934 | 0.0100 | 10.1 | 10.0 | 1.3 | 30.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.2869 | 0.2567 | 0.0100 | 17.9 | 20.0 | -10.5 | 30.0 |
| Benzo[a]anthracene | Ave | 1.055 | 0.999 | 0.8000 | 9.47 | 10.0 | -5.3 | 30.0 |
| Chrysene | Ave | 1.007 | 0.9502 | 0.7000 | 9.44 | 10.0 | -5.6 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-211933/11  Calibration Date: 12/22/2015 16:38

Instrument ID: A4HP9  Calib Start Date: 12/22/2015 12:48

GC Column: RXI-5SILMS  ID: 0.45(mm)  Calib End Date: 12/22/2015 20:02

Lab File ID: 51222011.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Ave | 1.620 | 1.613 | 0.0100 | 9.96 | 10.0 | -0.4 | 30.0 |
| Benzo[b]fluoranthene | Ave | 1.120 | 1.051 | 0.7000 | 9.38 | 10.0 | -6.2 | 30.0 |
| Benzo[k]fluoranthene | Ave | 1.192 | 1.120 | 0.7000 | 9.40 | 10.0 | -6.0 | 30.0 |
| Benzo[a]pyrene | Ave | 1.104 | 1.032 | 0.7000 | 9.35 | 10.0 | -6.5 | 30.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.102 | 1.016 | 0.5000 | 9.22 | 10.0 | -7.8 | 30.0 |
| Dibenz(a,h)anthracene | Ave | 0.8909 | 0.8702 | 0.4000 | 9.77 | 10.0 | -2.3 | 30.0 |
| Benzo[g,h,i]perylene | Ave | 0.9556 | 0.8941 | 0.5000 | 9.36 | 10.0 | -6.4 | 30.0 |
| 2-Fluorophenol (Surr) | Ave | 1.403 | 1.452 | | 10.3 | 10.0 | 3.5 | 30.0 |
| Phenol-d5 (Surr) | Ave | 1.824 | 1.913 | | 10.5 | 10.0 | 4.9 | 30.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.4835 | 0.4940 | | 10.2 | 10.0 | 2.2 | 30.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.216 | 1.204 | | 9.90 | 10.0 | -1.0 | 30.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1211 | 0.1301 | | 10.7 | 10.0 | 7.4 | 30.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7280 | 0.7281 | | 10.0 | 10.0 | 0.0 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212723/2 Calibration Date: 12/31/2015 09:36

Instrument ID: A4HP9 Calib Start Date: 12/22/2015 12:48

GC Column: RXI-5SILMS ID: 0.45(mm) Calib End Date: 12/22/2015 20:02

Lab File ID: 51231002.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.7095 | 0.5938 | | 8.37 | 10.0 | -16.3 | 20.0 |
| N-Nitrosodimethylamine | Ave | 1.048 | 0.9213 | | 8.79 | 10.0 | -12.1 | 20.0 |
| Pyridine | Ave | 1.874 | 1.700 | | 9.07 | 10.0 | -9.3 | 20.0 |
| Benzaldehyde | Ave | 1.371 | 1.205 | 0.0100 | 17.6 | 20.0 | -12.1 | 20.0 |
| Phenol | Ave | 2.021 | 1.865 | 0.8000 | 9.23 | 10.0 | -7.7 | 20.0 |
| Aniline | Ave | 2.494 | 2.350 | | 9.42 | 10.0 | -5.8 | 20.0 |
| Bis(2-chloroethyl)ether | Ave | 1.672 | 1.545 | 0.7000 | 9.24 | 10.0 | -7.6 | 20.0 |
| 2-Chlorophenol | Ave | 1.414 | 1.353 | 0.8000 | 9.57 | 10.0 | -4.3 | 20.0 |
| n-Decane | Ave | 1.810 | 1.710 | | 9.45 | 10.0 | -5.5 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.515 | 1.398 | | 9.22 | 10.0 | -7.8 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.507 | 1.411 | | 9.37 | 10.0 | -6.3 | 20.0 |
| Benzyl alcohol | Ave | 1.079 | 0.8990 | | 8.33 | 10.0 | -16.7 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.455 | 1.352 | | 9.29 | 10.0 | -7.1 | 20.0 |
| 2-Methylphenol | Ave | 1.471 | 1.383 | 0.7000 | 9.41 | 10.0 | -5.9 | 20.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.087 | 1.905 | | 9.13 | 10.0 | -8.7 | 20.0 |
| Indene | Ave | 2.527 | 2.437 | | 19.3 | 20.0 | -3.5 | 20.0 |
| 3 & 4 Methylphenol | Ave | 1.531 | 1.459 | | 9.53 | 10.0 | -4.7 | 20.0 |
| N-Nitrosodi-n-propylamine | Ave | 1.208 | 1.181 | 0.5000 | 9.78 | 10.0 | -2.2 | 20.0 |
| Acetophenone | Ave | 2.178 | 2.094 | 0.0100 | 9.61 | 10.0 | -3.9 | 20.0 |
| Hexachloroethane | Ave | 0.6335 | 0.6374 | 0.3000 | 10.1 | 10.0 | 0.6 | 20.0 |
| Nitrobenzene | Ave | 0.4609 | 0.4663 | 0.2000 | 10.1 | 10.0 | 1.2 | 20.0 |
| Isophorone | Ave | 0.8201 | 0.7964 | 0.4000 | 9.71 | 10.0 | -2.9 | 20.0 |
| 2-Nitrophenol | Ave | 0.1611 | 0.1649 | 0.1000 | 10.2 | 10.0 | 2.3 | 20.0 |
| 2,4-Dimethylphenol | Ave | 0.3844 | 0.3876 | 0.2000 | 10.1 | 10.0 | 0.8 | 20.0 |
| Benzoic acid | Ave | 0.2027 | 0.2112 | | 20.8 | 20.0 | 4.2 | 20.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4372 | 0.4292 | 0.3000 | 9.82 | 10.0 | -1.8 | 20.0 |
| 2,4-Dichlorophenol | Ave | 0.2571 | 0.2475 | 0.2000 | 9.63 | 10.0 | -3.7 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.2699 | 0.2493 | | 9.24 | 10.0 | -7.6 | 20.0 |
| Naphthalene | Ave | 1.070 | 0.995 | 0.7000 | 9.30 | 10.0 | -7.0 | 20.0 |
| 4-Chloroaniline | Ave | 0.4632 | 0.4391 | 0.0100 | 9.48 | 10.0 | -5.2 | 20.0 |
| 2,6-Dichlorophenol | Ave | 0.2521 | 0.2400 | | 9.52 | 10.0 | -4.8 | 20.0 |
| Hexachlorobutadiene | Ave | 0.1420 | 0.1394 | 0.0100 | 9.82 | 10.0 | -1.8 | 20.0 |
| Caprolactam | Ave | 0.1119 | 0.1073 | 0.0100 | 19.2 | 20.0 | -4.1 | 20.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3338 | 0.3389 | 0.2000 | 10.2 | 10.0 | 1.5 | 20.0 |
| 2-Methylnaphthalene | Ave | 0.6722 | 0.6325 | 0.4000 | 9.41 | 10.0 | -5.9 | 20.0 |
| 1-Methylnaphthalene | Ave | 0.6062 | 0.5495 | | 9.06 | 10.0 | -9.4 | 20.0 |
| Hexachlorocyclopentadiene | Ave | 0.2580 | 0.2835 | 0.0500 | 11.0 | 10.0 | 9.9 | 20.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4274 | 0.3944 | 0.0100 | 9.23 | 10.0 | -7.7 | 20.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3176 | 0.2974 | 0.2000 | 9.37 | 10.0 | -6.3 | 20.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3323 | 0.3121 | 0.2000 | 9.39 | 10.0 | -6.1 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212723/2 Calibration Date: 12/31/2015 09:36

Instrument ID: A4HP9 Calib Start Date: 12/22/2015 12:48

GC Column: RXI-5SILMS ID: 0.45(mm) Calib End Date: 12/22/2015 20:02

Lab File ID: 51231002.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.446 | 1.358 | 0.0100 | 9.39 | 10.0 | -6.1 | 20.0 |
| 2-Chloronaphthalene | Ave | 1.080 | 1.014 | 0.8000 | 9.39 | 10.0 | -6.1 | 20.0 |
| 2-Nitroaniline | Ave | 0.4431 | 0.4811 | 0.0100 | 10.9 | 10.0 | 8.6 | 20.0 |
| Dimethyl phthalate | Ave | 1.159 | 1.153 | 0.0100 | 9.95 | 10.0 | -0.5 | 20.0 |
| 1,3-Dinitrobenzene | Ave | 0.1847 | 0.1926 | | 10.4 | 10.0 | 4.2 | 20.0 |
| 2,6-Dinitrotoluene | Ave | 0.2660 | 0.2763 | | 10.4 | 10.0 | 3.9 | 20.0 |
| Acenaphthylene | Ave | 1.824 | 1.734 | 0.9000 | 9.51 | 10.0 | -4.9 | 20.0 |
| 3-Nitroaniline | Ave | 0.3099 | 0.3204 | 0.0100 | 10.3 | 10.0 | 3.4 | 20.0 |
| 2,4-Dinitrophenol | Lin1 | | 0.1585 | 0.0100 | 23.1 | 20.0 | 15.6 | 20.0 |
| Acenaphthene | Ave | 1.115 | 1.039 | 0.9000 | 9.32 | 10.0 | -6.8 | 20.0 |
| 4-Nitrophenol | Ave | 0.2041 | 0.2650 | | 26.0 | 20.0 | 29.8* | 20.0 |
| 2,4-Dinitrotoluene | Ave | 0.3379 | 0.3453 | 0.2000 | 10.2 | 10.0 | 2.2 | 20.0 |
| Dibenzofuran | Ave | 1.535 | 1.435 | 0.8000 | 9.35 | 10.0 | -6.5 | 20.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.2526 | 0.2442 | 0.0100 | 9.67 | 10.0 | -3.3 | 20.0 |
| Hexadecane | Ave | 0.9465 | 0.8857 | | 9.36 | 10.0 | -6.4 | 20.0 |
| Diethyl phthalate | Ave | 1.125 | 1.104 | 0.0100 | 9.82 | 10.0 | -1.8 | 20.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.5193 | 0.4776 | 0.4000 | 9.20 | 10.0 | -8.0 | 20.0 |
| 4-Nitroaniline | Ave | 0.2921 | 0.3404 | 0.0100 | 11.7 | 10.0 | 16.5 | 20.0 |
| Fluorene | Ave | 1.311 | 1.147 | 0.9000 | 8.75 | 10.0 | -12.5 | 20.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1203 | 0.0100 | 22.6 | 20.0 | 12.9 | 20.0 |
| N-Nitrosodiphenylamine | Ave | 0.5256 | 0.4920 | 0.0100 | 18.7 | 20.0 | -6.4 | 20.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 1.066 | 1.074 | | 10.1 | 10.0 | 0.8 | 20.0 |
| Azobenzene | Ave | 1.066 | 1.074 | | 10.1 | 10.0 | 0.8 | 20.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.1782 | 0.1710 | 0.1000 | 9.60 | 10.0 | -4.0 | 20.0 |
| Atrazine | Ave | 0.1673 | 0.1834 | 0.0100 | 21.9 | 20.0 | 9.6 | 20.0 |
| Hexachlorobenzene | Ave | 0.1903 | 0.1702 | 0.1000 | 8.95 | 10.0 | -10.5 | 20.0 |
| n-Octadecane | Ave | 0.6288 | 0.6282 | | 9.99 | 10.0 | -0.0 | 20.0 |
| Pentachlorophenol | Ave | 0.1138 | 0.1243 | 0.0500 | 21.8 | 20.0 | 9.2 | 20.0 |
| Phenanthrene | Ave | 1.082 | 1.003 | 0.7000 | 9.27 | 10.0 | -7.3 | 20.0 |
| Anthracene | Ave | 1.080 | 1.027 | 0.7000 | 9.51 | 10.0 | -4.9 | 20.0 |
| Carbazole | Ave | 0.9447 | 1.005 | 0.0100 | 10.6 | 10.0 | 6.4 | 20.0 |
| Di-n-butyl phthalate | Ave | 1.190 | 1.398 | 0.0100 | 11.7 | 10.0 | 17.5 | 20.0 |
| Fluoranthene | Ave | 1.067 | 1.067 | 0.6000 | 10.0 | 10.0 | -0.0 | 20.0 |
| Benzidine | Lin1 | | 0.5788 | | 21.7 | 20.0 | 8.5 | 20.0 |
| Pyrene | Ave | 1.056 | 1.074 | 0.6000 | 10.2 | 10.0 | 1.7 | 20.0 |
| Butyl benzyl phthalate | Ave | 0.5732 | 0.6229 | 0.0100 | 10.9 | 10.0 | 8.7 | 20.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.7835 | 0.8736 | 0.0100 | 11.1 | 10.0 | 11.5 | 20.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.2869 | 0.3245 | 0.0100 | 22.6 | 20.0 | 13.1 | 20.0 |
| Benzo[a]anthracene | Ave | 1.055 | 0.996 | 0.8000 | 9.44 | 10.0 | -5.6 | 20.0 |
| Chrysene | Ave | 1.007 | 0.9351 | 0.7000 | 9.29 | 10.0 | -7.1 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212723/2 | Calibration Date: 12/31/2015 09:36

Instrument ID: A4HP9 | Calib Start Date: 12/22/2015 12:48

GC Column: RXI-5SILMS ID: 0.45(mm) | Calib End Date: 12/22/2015 20:02

Lab File ID: 51231002.D | Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Ave | 1.620 | 1.730 | 0.0100 | 10.7 | 10.0 | 6.8 | 20.0 |
| Benzo[b]fluoranthene | Ave | 1.120 | 1.071 | 0.7000 | 9.56 | 10.0 | -4.4 | 20.0 |
| Benzo[k]fluoranthene | Ave | 1.192 | 1.091 | 0.7000 | 9.16 | 10.0 | -8.4 | 20.0 |
| Benzo[a]pyrene | Ave | 1.104 | 1.053 | 0.7000 | 9.54 | 10.0 | -4.6 | 20.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.102 | 1.144 | 0.5000 | 10.4 | 10.0 | 3.7 | 20.0 |
| Dibenz(a,h)anthracene | Ave | 0.8909 | 0.9461 | 0.4000 | 10.6 | 10.0 | 6.2 | 20.0 |
| Benzo[g,h,i]perylene | Ave | 0.9556 | 1.003 | 0.5000 | 10.5 | 10.0 | 4.9 | 20.0 |
| 2-Fluorophenol (Surr) | Ave | 1.403 | 1.332 | | 9.49 | 10.0 | -5.1 | 20.0 |
| Phenol-d5 (Surr) | Ave | 1.824 | 1.727 | | 9.47 | 10.0 | -5.3 | 20.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.4835 | 0.4840 | | 10.0 | 10.0 | 0.1 | 20.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.216 | 1.113 | | 9.15 | 10.0 | -8.5 | 20.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1211 | 0.1277 | | 10.5 | 10.0 | 5.4 | 20.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7280 | 0.7128 | | 9.79 | 10.0 | -2.1 | 20.0 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212548/7-A

Matrix: Solid | Lab File ID: 51231003.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30(g) | Date Analyzed: 12/31/2015 10:02

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 3.5 | U | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 9.5 | U | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 25 | U | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 8.9 | U | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 20 | U | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 20 | U | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 21 | U | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 17 | U | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 21 | U | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 0.45 | U | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 8.2 | U | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 0.50 | U | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 11 | U | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 9.1 | U | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 8.3 | U | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 18 | U | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 16 | U | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 9.2 | U | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 13 | U | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 21 | U | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 17 | U | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 13 | U | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 26 | U | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 17 | U | 330 | 17 |
| 83-32-9 | Acenaphthene | 0.76 | U | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 0.35 | U | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 9.2 | U | 100 | 9.2 |
| 120-12-7 | Anthracene | 0.78 | U | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 9.1 | U | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 12 | U | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 0.63 | U | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 0.64 | U | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 0.59 | U | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.35 | U | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212548/7-A

Matrix: Solid | Lab File ID: 51231003.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30(g) | Date Analyzed: 12/31/2015 10:02

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.68 | U | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 22 | U | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.0 | U | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 25.4 | J | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 10 | U | 70 | 10 |
| 105-60-2 | Caprolactam | 37 | U | 330 | 37 |
| 86-74-8 | Carbazole | 27 | U | 50 | 27 |
| 218-01-9 | Chrysene | 1.1 | U | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.66 | U | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 0.66 | U | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 16 | U | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 17 | U | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 15.9 | J | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 7.9 | U | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 0.55 | U | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 0.53 | U | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 2.1 | U | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 5.6 | U | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 8.1 | U | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 9.0 | U | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.35 | U | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 13.1 | J | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 6.3 | U | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 21 | U | 50 | 21 |
| 91-20-3 | Naphthalene | 0.82 | U | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 2.2 | U | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 9.1 | U | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 0.73 | U | 6.7 | 0.73 |
| 108-95-2 | Phenol | 7.3 | U | 50 | 7.3 |
| 129-00-0 | Pyrene | 0.44 | U | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 20 | U | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212548/7-A

Matrix: Solid | Lab File ID: 51231003.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30(g) | Date Analyzed: 12/31/2015 10:02

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 89 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 86 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 79 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 81 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 78 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 80 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212548/8-A

Matrix: Solid | Lab File ID: 51231004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30(g) | Date Analyzed: 12/31/2015 10:53

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 494 | | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 493 | | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 502 | | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 520 | | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 517 | | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 505 | | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 563 | | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 582 | | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 568 | | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 499 | | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 524 | | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 514 | | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 526 | | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 606 | | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 577 | | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 1100 | | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 536 | | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 1000 | | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 554 | | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 558 | | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 441 | | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 513 | | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 630 | | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 1450 | | 330 | 17 |
| 83-32-9 | Acenaphthene | 501 | | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 491 | | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 498 | | 100 | 9.2 |
| 120-12-7 | Anthracene | 528 | | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 1330 | | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 1090 | | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 537 | | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 522 | | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 524 | | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 567 | | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212548/8-A

Matrix: Solid | Lab File ID: 51231004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30(g) | Date Analyzed: 12/31/2015 10:53

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 525 | | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 546 | | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 501 | | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 642 | | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 609 | | 70 | 10 |
| 105-60-2 | Caprolactam | 1140 | | 330 | 37 |
| 86-74-8 | Carbazole | 638 | | 50 | 27 |
| 218-01-9 | Chrysene | 544 | | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 581 | | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 503 | | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 563 | | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 550 | | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 708 | | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 571 | | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 589 | | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 465 | | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 523 | | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 515 | | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 399 | | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 520 | | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 574 | | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 555 | | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 544 | | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 1100 | | 50 | 21 |
| 91-20-3 | Naphthalene | 501 | | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 545 | | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 1110 | | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 530 | | 6.7 | 0.73 |
| 108-95-2 | Phenol | 505 | | 50 | 7.3 |
| 129-00-0 | Pyrene | 574 | | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 516 | | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212548/8-A

Matrix: Solid | Lab File ID: 51231004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30(g) | Date Analyzed: 12/31/2015 10:53

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 81 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 80 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 82 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 77 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 74 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 90 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MS | Lab Sample ID: 240-59482-2 MS

Matrix: Solid | Lab File ID: 51231009.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:00

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 29.67(g) | Date Analyzed: 12/31/2015 12:58

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 631 | | 62 | 4.3 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 607 | | 120 | 12 |
| 95-95-4 | 2,4,5-Trichlorophenol | 627 | | 190 | 31 |
| 88-06-2 | 2,4,6-Trichlorophenol | 670 | | 190 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 640 | | 190 | 25 |
| 105-67-9 | 2,4-Dimethylphenol | 717 | | 190 | 25 |
| 51-28-5 | 2,4-Dinitrophenol | 1550 | | 410 | 26 |
| 121-14-2 | 2,4-Dinitrotoluene | 788 | | 250 | 21 |
| 606-20-2 | 2,6-Dinitrotoluene | 752 | | 250 | 26 |
| 91-58-7 | 2-Chloronaphthalene | 636 | | 62 | 0.56 |
| 95-57-8 | 2-Chlorophenol | 631 | | 62 | 10 |
| 91-57-6 | 2-Methylnaphthalene | 629 | | 8.3 | 0.62 |
| 95-48-7 | 2-Methylphenol | 623 | | 250 | 14 |
| 88-74-4 | 2-Nitroaniline | 808 | | 250 | 11 |
| 88-75-5 | 2-Nitrophenol | 710 | | 62 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 1320 | | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 490 | | 250 | 20 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 1620 | | 190 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 767 | | 62 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 727 | | 190 | 26 |
| 106-47-8 | 4-Chloroaniline | 396 | | 190 | 21 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 672 | | 62 | 16 |
| 100-01-6 | 4-Nitroaniline | 664 | | 250 | 32 |
| 100-02-7 | 4-Nitrophenol | 2010 | | 410 | 21 |
| 83-32-9 | Acenaphthene | 648 | | 8.3 | 0.94 |
| 208-96-8 | Acenaphthylene | 654 | | 8.3 | 0.43 |
| 98-86-2 | Acetophenone | 652 | | 120 | 11 |
| 120-12-7 | Anthracene | 711 | | 8.3 | 0.97 |
| 1912-24-9 | Atrazine | 1730 | | 250 | 11 |
| 100-52-7 | Benzaldehyde | 1330 | | 120 | 15 |
| 56-55-3 | Benzo[a]anthracene | 711 | | 8.3 | 0.78 |
| 50-32-8 | Benzo[a]pyrene | 713 | | 8.3 | 0.79 |
| 205-99-2 | Benzo[b]fluoranthene | 746 | | 8.3 | 0.73 |
| 191-24-2 | Benzo[g,h,i]perylene | 819 | | 8.3 | 0.43 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MS | Lab Sample ID: 240-59482-2 MS

Matrix: Solid | Lab File ID: 51231009.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:00

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 29.67(g) | Date Analyzed: 12/31/2015 12:58

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 712 | | 8.3 | 0.84 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 681 | | 120 | 27 |
| 111-44-4 | Bis(2-chloroethyl)ether | 697 | | 120 | 2.5 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 861 | | 87 | 24 |
| 85-68-7 | Butyl benzyl phthalate | 822 | | 87 | 12 |
| 105-60-2 | Caprolactam | 1490 | | 410 | 46 |
| 86-74-8 | Carbazole | 763 | | 62 | 33 |
| 218-01-9 | Chrysene | 717 | | 8.3 | 1.4 |
| 53-70-3 | Dibenz(a,h)anthracene | 844 | | 8.3 | 0.82 |
| 132-64-9 | Dibenzofuran | 669 | | 62 | 0.82 |
| 84-66-2 | Diethyl phthalate | 779 | | 87 | 20 |
| 131-11-3 | Dimethyl phthalate | 741 | | 87 | 21 |
| 84-74-2 | Di-n-butyl phthalate | 937 | | 87 | 19 |
| 117-84-0 | Di-n-octyl phthalate | 794 | | 87 | 9.8 |
| 206-44-0 | Fluoranthene | 781 | | 8.3 | 0.68 |
| 86-73-7 | Fluorene | 631 | | 8.3 | 0.66 |
| 118-74-1 | Hexachlorobenzene | 709 | | 8.3 | 2.6 |
| 87-68-3 | Hexachlorobutadiene | 629 | | 62 | 6.9 |
| 77-47-4 | Hexachlorocyclopentadiene | 333 | J | 410 | 10 |
| 67-72-1 | Hexachloroethane | 617 | | 62 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 811 | | 8.3 | 0.43 |
| 78-59-1 | Isophorone | 676 | | 62 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 674 | | 62 | 7.8 |
| 86-30-6 | N-Nitrosodiphenylamine | 1440 | | 62 | 26 |
| 91-20-3 | Naphthalene | 612 | | 8.3 | 1.0 |
| 98-95-3 | Nitrobenzene | 676 | | 120 | 2.7 |
| 87-86-5 | Pentachlorophenol | 1520 | | 190 | 11 |
| 85-01-8 | Phenanthrene | 715 | | 8.3 | 0.90 |
| 108-95-2 | Phenol | 601 | | 62 | 9.0 |
| 129-00-0 | Pyrene | 773 | | 8.3 | 0.55 |
| 15831-10-4 | 3 & 4 Methylphenol | 613 | | 500 | 25 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MS | Lab Sample ID: 240-59482-2 MS

Matrix: Solid | Lab File ID: 51231009.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:00

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 29.67(g) | Date Analyzed: 12/31/2015 12:58

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 89 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 80 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 84 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 78 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 77 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 101 | | 10-110 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MSD | Lab Sample ID: 240-59482-2 MSD

Matrix: Solid | Lab File ID: 51231010.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:00

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30.03(g) | Date Analyzed: 12/31/2015 13:23

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 636 | | 61 | 4.3 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 617 | | 120 | 12 |
| 95-95-4 | 2,4,5-Trichlorophenol | 643 | | 180 | 31 |
| 88-06-2 | 2,4,6-Trichlorophenol | 636 | | 180 | 11 |
| 120-83-2 | 2,4-Dichlorophenol | 646 | | 180 | 24 |
| 105-67-9 | 2,4-Dimethylphenol | 684 | | 180 | 24 |
| 51-28-5 | 2,4-Dinitrophenol | 1450 | | 400 | 26 |
| 121-14-2 | 2,4-Dinitrotoluene | 801 | | 240 | 21 |
| 606-20-2 | 2,6-Dinitrotoluene | 742 | | 240 | 26 |
| 91-58-7 | 2-Chloronaphthalene | 636 | | 61 | 0.55 |
| 95-57-8 | 2-Chlorophenol | 640 | | 61 | 10 |
| 91-57-6 | 2-Methylnaphthalene | 641 | | 8.2 | 0.61 |
| 95-48-7 | 2-Methylphenol | 638 | | 240 | 13 |
| 88-74-4 | 2-Nitroaniline | 791 | | 240 | 11 |
| 88-75-5 | 2-Nitrophenol | 752 | | 61 | 10 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 1470 | | 120 | 22 |
| 99-09-2 | 3-Nitroaniline | 589 | | 240 | 20 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 1590 | | 180 | 11 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 736 | | 61 | 16 |
| 59-50-7 | 4-Chloro-3-methylphenol | 726 | | 180 | 26 |
| 106-47-8 | 4-Chloroaniline | 378 | | 180 | 21 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 665 | | 61 | 16 |
| 100-01-6 | 4-Nitroaniline | 799 | | 240 | 32 |
| 100-02-7 | 4-Nitrophenol | 1740 | | 400 | 21 |
| 83-32-9 | Acenaphthene | 636 | | 8.2 | 0.93 |
| 208-96-8 | Acenaphthylene | 646 | | 8.2 | 0.43 |
| 98-86-2 | Acetophenone | 648 | | 120 | 11 |
| 120-12-7 | Anthracene | 704 | | 8.2 | 0.96 |
| 1912-24-9 | Atrazine | 1740 | | 240 | 11 |
| 100-52-7 | Benzaldehyde | 1270 | | 120 | 15 |
| 56-55-3 | Benzo[a]anthracene | 731 | | 8.2 | 0.77 |
| 50-32-8 | Benzo[a]pyrene | 714 | | 8.2 | 0.78 |
| 205-99-2 | Benzo[b]fluoranthene | 738 | | 8.2 | 0.72 |
| 191-24-2 | Benzo[g,h,i]perylene | 783 | | 8.2 | 0.43 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MSD | Lab Sample ID: 240-59482-2 MSD

Matrix: Solid | Lab File ID: 51231010.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:00

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30.03(g) | Date Analyzed: 12/31/2015 13:23

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 716 | | 8.2 | 0.83 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 683 | | 120 | 27 |
| 111-44-4 | Bis(2-chloroethyl)ether | 637 | | 120 | 2.4 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 938 | | 86 | 23 |
| 85-68-7 | Butyl benzyl phthalate | 832 | | 86 | 12 |
| 105-60-2 | Caprolactam | 1500 | | 400 | 45 |
| 86-74-8 | Carbazole | 828 | | 61 | 33 |
| 218-01-9 | Chrysene | 734 | | 8.2 | 1.3 |
| 53-70-3 | Dibenz(a,h)anthracene | 805 | | 8.2 | 0.81 |
| 132-64-9 | Dibenzofuran | 656 | | 61 | 0.81 |
| 84-66-2 | Diethyl phthalate | 764 | | 86 | 20 |
| 131-11-3 | Dimethyl phthalate | 730 | | 86 | 21 |
| 84-74-2 | Di-n-butyl phthalate | 939 | | 86 | 18 |
| 117-84-0 | Di-n-octyl phthalate | 788 | | 86 | 9.7 |
| 206-44-0 | Fluoranthene | 787 | | 8.2 | 0.67 |
| 86-73-7 | Fluorene | 619 | | 8.2 | 0.65 |
| 118-74-1 | Hexachlorobenzene | 682 | | 8.2 | 2.6 |
| 87-68-3 | Hexachlorobutadiene | 683 | | 61 | 6.9 |
| 77-47-4 | Hexachlorocyclopentadiene | 302 | J | 400 | 9.9 |
| 67-72-1 | Hexachloroethane | 646 | | 61 | 11 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 782 | | 8.2 | 0.43 |
| 78-59-1 | Isophorone | 691 | | 61 | 16 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 684 | | 61 | 7.7 |
| 86-30-6 | N-Nitrosodiphenylamine | 1420 | | 61 | 26 |
| 91-20-3 | Naphthalene | 623 | | 8.2 | 1.0 |
| 98-95-3 | Nitrobenzene | 680 | | 120 | 2.7 |
| 87-86-5 | Pentachlorophenol | 1280 | | 180 | 11 |
| 85-01-8 | Phenanthrene | 706 | | 8.2 | 0.89 |
| 108-95-2 | Phenol | 623 | | 61 | 8.9 |
| 129-00-0 | Pyrene | 766 | | 8.2 | 0.54 |
| 15831-10-4 | 3 & 4 Methylphenol | 618 | | 490 | 24 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MSD | Lab Sample ID: 240-59482-2 MSD

Matrix: Solid | Lab File ID: 51231010.D

Analysis Method: 8270D | Date Collected: 12/28/2015 10:00

Extract. Method: 3540C | Date Extracted: 12/30/2015 07:05

Sample wt/vol: 30.03(g) | Date Analyzed: 12/31/2015 13:23

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212723 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 89 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 79 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 86 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 77 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 78 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 90 | | 10-110 |

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A4HP9 Start Date: 12/22/2015 12:26

Analysis Batch Number: 211933 End Date: 12/22/2015 20:27

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| DFTPP 240-211933/1 | | 12/22/2015 12:26 | 1 | 51222001.D | RXI-5SILMS 0.45(mm) |
| STD5 240-211933/2 IC | | 12/22/2015 12:48 | 1 | 51222002.D | RXI-5SILMS 0.45(mm) |
| STD4 240-211933/3 IC | | 12/22/2015 13:13 | 1 | 51222003.D | RXI-5SILMS 0.45(mm) |
| STD3 240-211933/4 IC | | 12/22/2015 13:39 | 1 | 51222004.D | RXI-5SILMS 0.45(mm) |
| STD2 240-211933/5 IC | | 12/22/2015 14:04 | 1 | 51222005.D | RXI-5SILMS 0.45(mm) |
| STD1 240-211933/6 IC | | 12/22/2015 14:30 | 1 | 51222006.D | RXI-5SILMS 0.45(mm) |
| STD9 240-211933/7 IC | | 12/22/2015 14:56 | 1 | 51222007.D | RXI-5SILMS 0.45(mm) |
| STD8 240-211933/8 IC | | 12/22/2015 15:21 | 1 | 51222008.D | RXI-5SILMS 0.45(mm) |
| STD7 240-211933/9 IC | | 12/22/2015 15:47 | 1 | 51222009.D | RXI-5SILMS 0.45(mm) |
| STD6 240-211933/10 ICIS | | 12/22/2015 16:12 | 1 | 51222010.D | RXI-5SILMS 0.45(mm) |
| ICV 240-211933/11 | | 12/22/2015 16:38 | 1 | 51222011.D | RXI-5SILMS 0.45(mm) |
| STD5 240-211933/12 IC | | 12/22/2015 17:03 | 1 | | RXI-5SILMS 0.45(mm) |
| STD4 240-211933/13 IC | | 12/22/2015 17:29 | 1 | | RXI-5SILMS 0.45(mm) |
| STD3 240-211933/14 IC | | 12/22/2015 17:54 | 1 | | RXI-5SILMS 0.45(mm) |
| STD2 240-211933/15 IC | | 12/22/2015 18:20 | 1 | | RXI-5SILMS 0.45(mm) |
| STD6 240-211933/16 IC | | 12/22/2015 18:45 | 1 | | RXI-5SILMS 0.45(mm) |
| STD7 240-211933/17 IC | | 12/22/2015 19:10 | 1 | | RXI-5SILMS 0.45(mm) |
| STD8 240-211933/18 IC | | 12/22/2015 19:36 | 1 | | RXI-5SILMS 0.45(mm) |
| STD9 240-211933/19 IC | | 12/22/2015 20:02 | 1 | | RXI-5SILMS 0.45(mm) |
| ICV 240-211933/20 | | 12/22/2015 20:27 | 1 | | RXI-5SILMS 0.45(mm) |

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A4HP9 Start Date: 12/31/2015 09:15

Analysis Batch Number: 212723 End Date: 12/31/2015 17:11

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| DFTPP 240-212723/1 | | 12/31/2015 09:15 | 1 | 51231001.D | RXI-5SILMS 0.45(mm) |
| CCV 240-212723/2 CCVIS | | 12/31/2015 09:36 | 1 | 51231002.D | RXI-5SILMS 0.45(mm) |
| MB 240-212548/7-A | | 12/31/2015 10:02 | 1 | 51231003.D | RXI-5SILMS 0.45(mm) |
| LCS 240-212548/8-A | | 12/31/2015 10:53 | 1 | 51231004.D | RXI-5SILMS 0.45(mm) |
| 240-59482-3 | F-VB-UST-P-8 | 12/31/2015 12:08 | 1 | 51231007.D | RXI-5SILMS 0.45(mm) |
| 240-59482-2 | F-VB-UST-P-7 | 12/31/2015 12:33 | 1 | 51231008.D | RXI-5SILMS 0.45(mm) |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 12/31/2015 12:58 | 1 | 51231009.D | RXI-5SILMS 0.45(mm) |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 12/31/2015 13:23 | 1 | 51231010.D | RXI-5SILMS 0.45(mm) |
| ZZZZZ | | 12/31/2015 13:49 | 1 | | RXI-5SILMS 0.45(mm) |
| ZZZZZ | | 12/31/2015 14:14 | 1 | | RXI-5SILMS 0.45(mm) |
| ZZZZZ | | 12/31/2015 14:39 | 1 | | RXI-5SILMS 0.45(mm) |
| ZZZZZ | | 12/31/2015 15:04 | 1 | | RXI-5SILMS 0.45(mm) |
| ZZZZZ | | 12/31/2015 15:30 | 1 | | RXI-5SILMS 0.45(mm) |
| ZZZZZ | | 12/31/2015 15:55 | 1 | | RXI-5SILMS 0.45(mm) |
| ZZZZZ | | 12/31/2015 16:20 | 1 | | RXI-5SILMS 0.45(mm) |
| ZZZZZ | | 12/31/2015 16:45 | 1 | | RXI-5SILMS 0.45(mm) |
| ZZZZZ | | 12/31/2015 17:11 | 1 | | RXI-5SILMS 0.45(mm) |

# GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Batch Number: 212548  Batch Start Date: 12/30/15 08:55  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 12/31/15 12:55

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exBENZALDEHYD 00033 | exBNASPIKE 00055 | exBNASURR W 00030 | |
|---|---|---|---|---|---|---|---|---|---|
| 240-59482-2 | F-VB-UST-P-7 | 3540C, 8270D | T | 30.06 g | 2 mL | | | 1 mL | |
| 240-59482-3 | F-VB-UST-P-8 | 3540C, 8270D | T | 30.35 g | 2 mL | | | 1 mL | |
| 240-59482-2 MS | F-VB-UST-P-7 | 3540C, 8270D | T | 29.67 g | 2 mL | 1 mL | 1 mL | 1 mL | |
| 240-59482-2 MSD | F-VB-UST-P-7 | 3540C, 8270D | T | 30.03 g | 2 mL | 1 mL | 1 mL | 1 mL | |
| MB 240-212548/7 | | 3540C, 8270D | | 30 g | 2 mL | | | 1 mL | |
| LCS 240-212548/8 | | 3540C, 8270D | | 30 g | 2 mL | 1 mL | 1 mL | 1 mL | |

| Batch Notes | |
|---|---|
| Balance ID | 1339911 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2348898 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8015C - GRO

## Nonhalogenated Organics using GC/FID - Modified (Gasoline Range Organics) by Method 8015C

FORM II
GC VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): RTX-VRX ID: 0.45(mm)

| Client Sample ID | Lab Sample ID | TFT1 # |
|---|---|---|
| F-VB-UST-P-7 | 240-59482-2 | 70 |
| F-VB-UST-P-8 | 240-59482-3 | 67 |
| | MB 240-212559/1-A | 68 |
| | LCS 240-212559/2-A | 90 |
| F-VB-UST-P-7 MS | 240-59482-2 MS | 82 |
| F-VB-UST-P-7 MSD | 240-59482-2 MSD | 82 |

| | QC LIMITS |
|---|---|
| TFT = Trifluorotoluene (Surr) | 40-139 |

# Column to be used to flag recovery values

FORM II 8015C

FORM III
GC VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: Y0122920.D

Lab ID: LCS 240-212559/2-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 771 | 96 | 74-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: Y0122922.D

Lab ID: 240-59482-2 MS Client ID: F-VB-UST-P-7 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 979 | 56 U | 885 | 90 | 10-136 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: Y0122923.D

Lab ID: 240-59482-2 MSD Client ID: F-VB-UST-P-7 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 973 | 841 | 86 | 5 | 40 | 10-136 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM IV
GC VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: MB 240-212559/1-A

Matrix: Solid  Date Extracted: 12/30/2015 07:39

Lab File ID:(1) Y0122919.D  Lab File ID:(2)

Date Analyzed:(1) 12/30/2015 09:18  Date Analyzed:(2)

Instrument ID:(1) YPID  Instrument ID:(2)

GC Column:(1) RTX-VRX  ID: 0.45(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-212559/2-A | 12/30/2015 09:50 | |
| F-VB-UST-P-7 | 240-59482-2 | 12/30/2015 10:23 | |
| F-VB-UST-P-7 MS | 240-59482-2 MS | 12/30/2015 10:57 | |
| F-VB-UST-P-7 MSD | 240-59482-2 MSD | 12/30/2015 11:31 | |
| F-VB-UST-P-8 | 240-59482-3 | 12/30/2015 12:04 | |

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Sample No.: CCVRT 240-212533/4  Date Analyzed: 12/30/2015 06:31

Instrument ID: YPID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): Y0122914.D  Heated Purge: (Y/N) N

Calibration ID: 31080

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 9.73 | | |
| UPPER LIMIT | | | | 9.83 | | |
| LOWER LIMIT | | | | 9.63 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCVRT 240-212533/4 | | 12/30/2015 06:31 | Y0122914.D | 9.73 | | |
| CCV 240-212533/8 | | 12/30/2015 08:44 | Y0122918.D | 9.74 | | |
| MB 240-212559/1-A | | 12/30/2015 09:18 | Y0122919.D | 9.73 | | |
| LCS 240-212559/2-A | | 12/30/2015 09:50 | Y0122920.D | 9.73 | | |
| 240-59482-2 | F-VB-UST-P-7 | 12/30/2015 10:23 | Y0122921.D | 9.73 | | |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 12/30/2015 10:57 | Y0122922.D | 9.74 | | |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 12/30/2015 11:31 | Y0122923.D | 9.73 | | |
| 240-59482-3 | F-VB-UST-P-8 | 12/30/2015 12:04 | Y0122924.D | 9.73 | | |
| CCV 240-212533/17 | | 12/30/2015 13:44 | Y0122927.D | 9.74 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 | Lab Sample ID: 240-59482-2

Matrix: Solid | Lab File ID: Y0122921.D

Analysis Method: 8015C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.05(g) | Date Analyzed: 12/30/2015 10:23

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212533 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 56 | U | 120 | 56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 70 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-8 | Lab Sample ID: 240-59482-3

Matrix: Solid | Lab File ID: Y0122924.D

Analysis Method: 8015C | Date Collected: 12/28/2015 10:30

Sample wt/vol: 5.00(g) | Date Analyzed: 12/30/2015 12:04

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 16.1 | Level: (low/med) Low

Analysis Batch No.: 212533 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 67 | | 40-139 |

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 201382

SDG No.:

Instrument ID: YPID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/12/2015 10:39 Calibration End Date: 10/12/2015 13:35 Calibration ID: 31080

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-201382/4 | Y0101204.D |
| Level 2 | STD2 240-201382/5 | Y0101205.D |
| Level 3 | STD3 240-201382/6 | Y0101206.D |
| Level 4 | STD4 240-201382/7 | Y0101207.D |
| Level 5 | STD5 240-201382/8 | Y0101208.D |
| Level 6 | STD6 240-201382/9 | Y0101209.D |

| ANALYTE | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 6 | | | | | RT WINDOW | AVG RT |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 12.032 | 12.032 | 12.032 | 12.032 | 12.032 | 12.032 | | | | | 4.928 - 19.136 | 12.032 |
| C6-C12 | 13.458 | 13.458 | 13.458 | 13.458 | 13.458 | 13.458 | | | | | 4.928 - 21.988 | 13.458 |
| Trifluorotoluene (Surr) | 9.720 | 9.722 | 9.720 | 9.726 | 9.719 | 9.722 | | | | | 9.476 - 9.976 | 9.722 |

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 201382

SDG No.:

Instrument ID: YPID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/12/2015 10:39 Calibration End Date: 10/12/2015 13:35 Calibration ID: 31080

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-201382/4 | Y0101204.D |
| Level 2 | STD2 240-201382/5 | Y0101205.D |
| Level 3 | STD3 240-201382/6 | Y0101206.D |
| Level 4 | STD4 240-201382/7 | Y0101207.D |
| Level 5 | STD5 240-201382/8 | Y0101208.D |
| Level 6 | STD6 240-201382/9 | Y0101209.D |

| ANALYTE | CF | | | | CURVE TYPE | COEFFICIENT | | | # | MIN CF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 5 | LVL 2<br>LVL 6 | LVL 3 | LVL 4 | | B | M1 | M2 | | | | | | | | |
| Gasoline Range Organics [C6 - C10] | 188416<br>173575 | 165916<br>171151 | 176757 | 171370 | Ave | | 174530.549 | | | | 4.4 | | 20.0 | | | |
| C6-C12 | 222173<br>187083 | 182916<br>184022 | 192254 | 185117 | Ave | | 192260.693 | | | | 7.8 | | 20.0 | | | |
| Trifluorotoluene (Surr) | 210718<br>163918 | 183500<br>159255 | 169226 | 148916 | Ave | | 172588.858 | | | | 12.7 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 201382

SDG No.:

Instrument ID: YPID GC Column: RTX-VRX ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/12/2015 10:39 Calibration End Date: 10/12/2015 13:35 Calibration ID: 31080

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-201382/4 | Y0101204.D |
| Level 2 | STD2 240-201382/5 | Y0101205.D |
| Level 3 | STD3 240-201382/6 | Y0101206.D |
| Level 4 | STD4 240-201382/7 | Y0101207.D |
| Level 5 | STD5 240-201382/8 | Y0101208.D |
| Level 6 | STD6 240-201382/9 | Y0101209.D |

| ANALYTE | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (UG/L) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|
| | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Gasoline Range Organics [C6 - C10] | Ave | 18841554<br>273840999 | 33183190 | 70702651 | 137095852 | 208289684 | 100<br>1600 | 200 | 400 | 800 | 1200 |
| C6-C12 | Ave | 22217288<br>294434655 | 36583184 | 76901726 | 148093223 | 224499426 | 100<br>1600 | 200 | 400 | 800 | 1200 |
| Trifluorotoluene (Surr) | Ave | 210718<br>7962770 | 367000 | 846131 | 1489159 | 3278353 | 1.00<br>50.0 | 2.00 | 5.00 | 10.0 | 20.0 |

Curve Type Legend:
Ave = Average

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-201382/10 Calibration Date: 10/12/2015 14:09

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0101210.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 174531 | 167422 | | 767 | 800 | -4.1 | 20.0 |
| C6-C12 | Ave | 192261 | 181216 | | 754 | 800 | -5.7 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 172589 | 166182 | | 9.63 | 10.0 | -3.7 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-201382/10 Calibration Date: 10/12/2015 14:09

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0101210.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 12.03 | 4.93 | 19.14 |
| C6-C12 | 13.46 | 4.93 | 21.99 |
| Trifluorotoluene (Surr) | 9.72 | 9.48 | 9.98 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCVRT 240-212533/4 Calibration Date: 12/30/2015 06:31

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0122914.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 174531 | 177649 | | 814 | 800 | 1.8 | 20.0 |
| C6-C12 | Ave | 192261 | 193248 | | 804 | 800 | 0.5 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 172589 | 139370 | | 8.08 | 10.0 | -19.2 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCVRT 240-212533/4 Calibration Date: 12/30/2015 06:31

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0122914.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 12.01 | 4.90 | 19.12 |
| C6-C12 | 13.44 | 4.90 | 21.98 |
| Trifluorotoluene (Surr) | 9.73 | 9.48 | 9.98 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212533/8 Calibration Date: 12/30/2015 08:44

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0122918.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 174531 | 179834 | | 824 | 800 | 3.0 | 20.0 |
| C6-C12 | Ave | 192261 | 193369 | | 805 | 800 | 0.6 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 172589 | 138663 | | 8.03 | 10.0 | -19.7 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212533/8 Calibration Date: 12/30/2015 08:44

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0122918.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 12.01 | 4.90 | 19.12 |
| C6-C12 | 13.44 | 4.90 | 21.98 |
| Trifluorotoluene (Surr) | 9.74 | 9.48 | 9.98 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212533/17 Calibration Date: 12/30/2015 13:44

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0122927.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 174531 | 159489 | | 731 | 800 | -8.6 | 20.0 |
| C6-C12 | Ave | 192261 | 172448 | | 718 | 800 | -10.3 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 172589 | 139318 | | 8.07 | 10.0 | -19.3 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212533/17 Calibration Date: 12/30/2015 13:44

Instrument ID: YPID Calib Start Date: 10/12/2015 10:39

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 10/12/2015 13:35

Lab File ID: Y0122927.D Heated Purge: (Y/N) N

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Gasoline Range Organics [C6 - C10] | 12.01 | 4.90 | 19.12 |
| C6-C12 | 13.44 | 4.90 | 21.98 |
| Trifluorotoluene (Surr) | 9.74 | 9.48 | 9.98 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-212559/1-A

Matrix: Solid  Lab File ID: Y0122919.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 12/30/2015 09:18

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 212533  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 46 | U | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 68 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-212559/2-A

Matrix: Solid  Lab File ID: Y0122920.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 12/30/2015 09:50

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 212533  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 771 | | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 90 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MS | Lab Sample ID: 240-59482-2 MS

Matrix: Solid | Lab File ID: Y0122922.D

Analysis Method: 8015C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.01(g) | Date Analyzed: 12/30/2015 10:57

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212533 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 885 | | 120 | 56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 82 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 MSD | Lab Sample ID: 240-59482-2 MSD

Matrix: Solid | Lab File ID: Y0122923.D

Analysis Method: 8015C | Date Collected: 12/28/2015 10:00

Sample wt/vol: 5.04(g) | Date Analyzed: 12/30/2015 11:31

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 18.4 | Level: (low/med) Low

Analysis Batch No.: 212533 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 841 | | 120 | 56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 82 | | 40-139 |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: YPID Start Date: 10/12/2015 10:39

Analysis Batch Number: 201382 End Date: 10/12/2015 15:19

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-201382/4 IC | | 10/12/2015 10:39 | 1 | Y0101204.D | RTX-VRX 0.45(mm) |
| STD2 240-201382/5 IC | | 10/12/2015 11:13 | 1 | Y0101205.D | RTX-VRX 0.45(mm) |
| STD3 240-201382/6 IC | | 10/12/2015 11:49 | 1 | Y0101206.D | RTX-VRX 0.45(mm) |
| STD4 240-201382/7 IC | | 10/12/2015 12:23 | 1 | Y0101207.D | RTX-VRX 0.45(mm) |
| STD5 240-201382/8 IC | | 10/12/2015 12:59 | 1 | Y0101208.D | RTX-VRX 0.45(mm) |
| STD6 240-201382/9 IC | | 10/12/2015 13:35 | 1 | Y0101209.D | RTX-VRX 0.45(mm) |
| ICV 240-201382/10 | | 10/12/2015 14:09 | 1 | Y0101210.D | RTX-VRX 0.45(mm) |
| ZZZZZ | | 10/12/2015 14:45 | 2 | | RTX-VRX 0.45(mm) |
| CCV 240-201382/12 | | 10/12/2015 15:19 | 1 | | RTX-VRX 0.45(mm) |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: YPID Start Date: 12/30/2015 06:31

Analysis Batch Number: 212533 End Date: 12/30/2015 13:44

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCVRT 240-212533/4 | | 12/30/2015 06:31 | 1 | Y0122914.D | RTX-VRX 0.45(mm) |
| CCV 240-212533/8 | | 12/30/2015 08:44 | 1 | Y0122918.D | RTX-VRX 0.45(mm) |
| MB 240-212559/1-A | | 12/30/2015 09:18 | 1 | Y0122919.D | RTX-VRX 0.45(mm) |
| LCS 240-212559/2-A | | 12/30/2015 09:50 | 1 | Y0122920.D | RTX-VRX 0.45(mm) |
| 240-59482-2 | F-VB-UST-P-7 | 12/30/2015 10:23 | 1 | Y0122921.D | RTX-VRX 0.45(mm) |
| 240-59482-2 MS | F-VB-UST-P-7 MS | 12/30/2015 10:57 | 1 | Y0122922.D | RTX-VRX 0.45(mm) |
| 240-59482-2 MSD | F-VB-UST-P-7 MSD | 12/30/2015 11:31 | 1 | Y0122923.D | RTX-VRX 0.45(mm) |
| 240-59482-3 | F-VB-UST-P-8 | 12/30/2015 12:04 | 1 | Y0122924.D | RTX-VRX 0.45(mm) |
| ZZZZZ | | 12/30/2015 13:10 | 1 | | RTX-VRX 0.45(mm) |
| CCV 240-212533/17 | | 12/30/2015 13:44 | 1 | Y0122927.D | RTX-VRX 0.45(mm) |

# GC VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Batch Number: 212559  Batch Start Date: 12/30/15 07:39  Batch Analyst: Reed, Rachel T

Batch Method: 5030A  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| MB 240-212559/1 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| LCS 240-212559/2 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| 240-59482-2 | F-VB-UST-P-7 | 5030A, 8015C | T | 5.05 g | 5 mL | | | | |
| 240-59482-2 MS | F-VB-UST-P-7 | 5030A, 8015C | T | 5.01 g | 5 mL | | | | |
| 240-59482-2 MSD | F-VB-UST-P-7 | 5030A, 8015C | T | 5.04 g | 5 mL | | | | |
| 240-59482-3 | F-VB-UST-P-8 | 5030A, 8015C | T | 5.00 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| Balance ID | B045 |
| Blank Soil Lot Number | 1757979 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8015C - DRO

## Nonhalogenated Organics using GC/FID - Modified (Diesel Range Organics) by Method 8015C

FORM II
DIESEL RANGE ORGANICS SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): ID:

| Client Sample ID | Lab Sample ID | OTPH1 # |
|---|---|---|
| F-VB-UST-P-7 | 240-59482-2 | 92 |
| F-VB-UST-P-8 | 240-59482-3 | 104 |
| | MB 240-212549/8-A | 90 |
| | LCS 240-212549/9-A | 97 |

| | QC LIMITS |
|---|---|
| OTPH = o-Terphenyl (Surr) | 40-160 |

# Column to be used to flag recovery values

FORM II 8015C

FORM III
DIESEL RANGE ORGANICS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5F23116.D

Lab ID: LCS 240-212549/9-A Client ID:

| COMPOUND | SPIKE ADDED (mg/Kg) | LCS CONCENTRATION (mg/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Diesel | 84.0 | 75.4 | 90 | 52-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM IV
DIESEL RANGE ORGANICS METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: MB 240-212549/8-A

Matrix: Solid Date Extracted: 12/30/2015 07:14

Lab File ID:(1) P5F23115.D Lab File ID:(2)

Date Analyzed:(1) 12/31/2015 20:19 Date Analyzed:(2)

Instrument ID:(1) A2HP5F Instrument ID:(2)

GC Column:(1) ID: GC Column:(2) ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| F-VB-UST-P-7 | 240-59482-2 | 12/31/2015 17:44 | |
| F-VB-UST-P-8 | 240-59482-3 | 12/31/2015 18:15 | |
| | LCS 240-212549/9-A | 12/31/2015 20:50 | |

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Sample No.: STD3 240-210457/10 Date Analyzed: 12/11/2015 18:19

Instrument ID: A2HP5F GC Column: ID: ()

Lab File ID (Standard): P5F21110.D Heated Purge: (Y/N) N

Calibration ID: 32080

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| INITIAL CALIBRATION SURROGATE | | | | 9.84 | | |
| UPPER LIMIT | | | | 9.94 | | |
| LOWER LIMIT | | | | 9.74 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| STD3 240-210457/10 ICRT | | 12/11/2015 18:19 | P5F21110.D | 9.84 | | |

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Sample No.: CCV 240-212780/7  Date Analyzed: 12/31/2015 16:11

Instrument ID: A2HP5F  GC Column:  ID: ()

Lab File ID (Standard): P5F23107.D  Heated Purge: (Y/N) N

Calibration ID: 32080

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCV 240-212780/7 CCVRT | | 12/31/2015 16:11 | P5F23107.D | 9.83 | | |
| 240-59482-2 | F-VB-UST-P-7 | 12/31/2015 17:44 | P5F23110.D | 9.83 | | |
| 240-59482-3 | F-VB-UST-P-8 | 12/31/2015 18:15 | P5F23111.D | 9.83 | | |
| MB 240-212549/8-A | | 12/31/2015 20:19 | P5F23115.D | 9.83 | | |
| LCS 240-212549/9-A | | 12/31/2015 20:50 | P5F23116.D | 9.83 | | |
| CCV 240-212780/17 | | 12/31/2015 21:21 | P5F23117.D | 9.83 | | |

OTPH = o-Terphenyl

OTPH RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 | Lab Sample ID: 240-59482-2

Matrix: Solid | Lab File ID: P5F23110.D

Analysis Method: 8015C | Date Collected: 12/28/2015 10:00

Extraction Method: 3540C | Date Extracted: 12/30/2015 07:14

Sample wt/vol: 29.97(g) | Date Analyzed: 12/31/2015 17:44

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212780 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.9 | U | 20 | 7.9 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 92 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-8 | Lab Sample ID: 240-59482-3

Matrix: Solid | Lab File ID: P5F23111.D

Analysis Method: 8015C | Date Collected: 12/28/2015 10:30

Extraction Method: 3540C | Date Extracted: 12/30/2015 07:14

Sample wt/vol: 29.78(g) | Date Analyzed: 12/31/2015 18:15

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212780 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 42 | | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 104 | | 40-160 |

FORM VI
DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210457

SDG No.:

Instrument ID: A2HP5F GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 12/11/2015 17:00 Calibration End Date: 12/11/2015 19:39 Calibration ID: 32080

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-210457/8 | P5F21108.D |
| Level 2 | STD2 240-210457/9 | P5F21109.D |
| Level 3 | STD3 240-210457/10 | P5F21110.D |
| Level 4 | STD4 240-210457/11 | P5F21111.D |
| Level 5 | STD5 240-210457/12 | P5F21112.D |

| ANALYTE | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | | | | | RT WINDOW | AVG RT |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | 9.334 | 9.334 | 9.334 | 9.334 | 9.334 | | | | | | 5.118 - 13.551 | 9.334 |
| o-Terphenyl (Surr) | 9.835 | 9.836 | 9.837 | 9.838 | 9.840 | | | | | | 9.735 - 9.935 | 9.837 |

FORM VI
DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210457

SDG No.:

Instrument ID: A2HP5F GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 12/11/2015 17:00 Calibration End Date: 12/11/2015 19:39 Calibration ID: 32080

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-210457/8 | P5F21108.D |
| Level 2 | STD2 240-210457/9 | P5F21109.D |
| Level 3 | STD3 240-210457/10 | P5F21110.D |
| Level 4 | STD4 240-210457/11 | P5F21111.D |
| Level 5 | STD5 240-210457/12 | P5F21112.D |

| ANALYTE | CF | | | | CURVE TYPE | COEFFICIENT | | | # | MIN CF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 5 | LVL 2 | LVL 3 | LVL 4 | | B | M1 | M2 | | | | | | | | |
| Diesel Range Organics [C10 - C28] | 156544 134522 | 139315 | 139028 | 130044 | Ave | | 139890.741 | | | | 7.2 | | 20.0 | | | |
| o-Terphenyl (Surr) | 166040 172555 | 155445 | 168395 | 159963 | Ave | | 164479.356 | | | | 4.1 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 210457

SDG No.:

Instrument ID: A2HP5F GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 12/11/2015 17:00 Calibration End Date: 12/11/2015 19:39 Calibration ID: 32080

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-210457/8 | P5F21108.D |
| Level 2 | STD2 240-210457/9 | P5F21109.D |
| Level 3 | STD3 240-210457/10 | P5F21110.D |
| Level 4 | STD4 240-210457/11 | P5F21111.D |
| Level 5 | STD5 240-210457/12 | P5F21112.D |

| ANALYTE | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|
| | | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| Diesel Range Organics [C10 - C28] | Ave | 15654411 | 27863003 | 69514146 | 130044202 | 269044173 | 100 | 200 | 500 | 1000 | 2000 |
| o-Terphenyl (Surr) | Ave | 1328317 | 2487115 | 6735786 | 12797061 | 27608729 | 8.00 | 16.0 | 40.0 | 80.0 | 160 |

Curve Type Legend:
Ave = Average

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-210457/13  Calibration Date: 12/11/2015 20:19

Instrument ID: A2HP5F  Calib Start Date: 12/11/2015 17:00

GC Column:  ID:  Calib End Date: 12/11/2015 19:39

Lab File ID: P5F21113.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 139891 | 132891 | | 476 | 502 | -5.0 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-210457/13 Calibration Date: 12/11/2015 20:19

Instrument ID: A2HP5F Calib Start Date: 12/11/2015 17:00

GC Column: ID: Calib End Date: 12/11/2015 19:39

Lab File ID: P5F21113.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.33 | 5.12 | 13.55 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212780/7  Calibration Date: 12/31/2015 16:11

Instrument ID: A2HP5F  Calib Start Date: 12/11/2015 17:00

GC Column:  ID:  Calib End Date: 12/11/2015 19:39

Lab File ID: P5F23107.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 139891 | 142174 | | 508 | 500 | 1.6 | 20.0 |
| o-Terphenyl (Surr) | Ave | 164479 | 176161 | | 42.8 | 40.0 | 7.1 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212780/7 Calibration Date: 12/31/2015 16:11

Instrument ID: A2HP5F Calib Start Date: 12/11/2015 17:00

GC Column: ID: Calib End Date: 12/11/2015 19:39

Lab File ID: P5F23107.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.32 | 5.10 | 13.54 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212780/17 Calibration Date: 12/31/2015 21:21

Instrument ID: A2HP5F Calib Start Date: 12/11/2015 17:00

GC Column: ID: Calib End Date: 12/11/2015 19:39

Lab File ID: P5F23117.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 139891 | 144589 | | 517 | 500 | 3.4 | 20.0 |
| o-Terphenyl (Surr) | Ave | 164479 | 186811 | | 45.4 | 40.0 | 13.6 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212780/17 Calibration Date: 12/31/2015 21:21

Instrument ID: A2HP5F Calib Start Date: 12/11/2015 17:00

GC Column: ID: Calib End Date: 12/11/2015 19:39

Lab File ID: P5F23117.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| Diesel Range Organics [C10 - C28] | 9.32 | 5.10 | 13.54 |
| o-Terphenyl (Surr) | 9.83 | 9.73 | 9.93 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212549/8-A

Matrix: Solid | Lab File ID: P5F23115.D

Analysis Method: 8015C | Date Collected:

Extraction Method: 3540C | Date Extracted: 12/30/2015 07:14

Sample wt/vol: 30(g) | Date Analyzed: 12/31/2015 20:19

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212780 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 6.4 | U | 17 | 6.4 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 90 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212549/9-A

Matrix: Solid | Lab File ID: P5F23116.D

Analysis Method: 8015C | Date Collected:

Extraction Method: 3540C | Date Extracted: 12/30/2015 07:14

Sample wt/vol: 30(g) | Date Analyzed: 12/31/2015 20:50

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212780 | Units: mg/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 97 | | 40-160 |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A2HP5F Start Date: 12/11/2015 17:00

Analysis Batch Number: 210457 End Date: 12/12/2015 02:17

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-210457/8 IC | | 12/11/2015 17:00 | 1 | P5F21108.D | |
| STD2 240-210457/9 IC | | 12/11/2015 17:40 | 1 | P5F21109.D | |
| STD3 240-210457/10 ICRT | | 12/11/2015 18:19 | 1 | P5F21110.D | |
| STD4 240-210457/11 IC | | 12/11/2015 18:59 | 1 | P5F21111.D | |
| STD5 240-210457/12 IC | | 12/11/2015 19:39 | 1 | P5F21112.D | |
| ICV 240-210457/13 | | 12/11/2015 20:19 | 1 | P5F21113.D | |
| ZZZZZ | | 12/11/2015 20:58 | 1 | | |
| ZZZZZ | | 12/11/2015 21:38 | 1 | | |
| ZZZZZ | | 12/11/2015 23:37 | 1 | | |
| ZZZZZ | | 12/12/2015 00:17 | 1 | | |
| ZZZZZ | | 12/12/2015 00:57 | 1 | | |
| ZZZZZ | | 12/12/2015 01:37 | 1 | | |
| CCV 240-210457/22 | | 12/12/2015 02:17 | 1 | | |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A2HP5F Start Date: 12/31/2015 16:11

Analysis Batch Number: 212780 End Date: 12/31/2015 21:21

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-212780/7 CCVRT | | 12/31/2015 16:11 | 1 | P5F23107.D | |
| ZZZZZ | | 12/31/2015 16:42 | 10 | | |
| ZZZZZ | | 12/31/2015 17:13 | 1 | | |
| 240-59482-2 | F-VB-UST-P-7 | 12/31/2015 17:44 | 1 | P5F23110.D | |
| 240-59482-3 | F-VB-UST-P-8 | 12/31/2015 18:15 | 1 | P5F23111.D | |
| ZZZZZ | | 12/31/2015 18:46 | 1 | | |
| ZZZZZ | | 12/31/2015 19:17 | 10 | | |
| ZZZZZ | | 12/31/2015 19:47 | 10 | | |
| MB 240-212549/8-A | | 12/31/2015 20:19 | 1 | P5F23115.D | |
| LCS 240-212549/9-A | | 12/31/2015 20:50 | 1 | P5F23116.D | |
| CCV 240-212780/17 | | 12/31/2015 21:21 | 1 | P5F23117.D | |

# DIESEL RANGE ORGANICS BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Batch Number: 212549  Batch Start Date: 12/30/15 08:55  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 12/31/15 12:55

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exDIESELSPKw 00069 | exOTPSURRW 00007 | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-59482-2 | F-VB-UST-P-7 | 3540C, 8015C | T | 29.97 g | 5 mL | | 1 mL | | |
| 240-59482-3 | F-VB-UST-P-8 | 3540C, 8015C | T | 29.78 g | 5 mL | | 1 mL | | |
| MB 240-212549/8 | | 3540C, 8015C | | 30 g | 5 mL | | 1 mL | | |
| LCS 240-212549/9 | | 3540C, 8015C | | 30 g | 5 mL | 1 mL | 1 mL | | |

| Batch Notes | |
|---|---|
| Balance ID | 1339911 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2348898 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Method 8082A

## Polychlorinated Biphenyls (PCBs) (GC) by Method 8082A

FORM II
PCBS SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low

GC Column (2): CLP-1 ID: 0.53(mm)

| Client Sample ID | Lab Sample ID | TCX2 # | DCB2 # |
|---|---|---|---|
| F-VB-UST-P-7 | 240-59482-2 | 84 | 72 |
| F-VB-UST-P-8 | 240-59482-3 | 57 | 92 |
| | MB 240-212544/7-A | 106 | 110 |
| | LCS 240-212544/8-A | 102 | 110 |
| | 240-59503-G-4-B MS | 44 | 43 |
| | 240-59503-G-4-C MSD | 31 p | 37 |

| | QC LIMITS |
|---|---|
| TCX = Tetrachloro-m-xylene | 13-134 |
| DCB = DCB Decachlorobiphenyl | 10-155 |

# Column to be used to flag recovery values

FORM II 8082A

FORM III
PCBS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P1200020.D

Lab ID: LCS 240-212544/8-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Aroclor-1016 | 333 | 314 | 94 | 51-120 | |
| Aroclor-1260 | 333 | 337 | 101 | 48-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM III
PCBS MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P1200017.D

Lab ID: 240-59503-G-4-B MS Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Aroclor-1016 | 442 | 16 U | 175 | 40 | 17-141 | |
| Aroclor-1260 | 442 | 12 U | 143 | 32 | 19-137 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM III
PCBS MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P1200018.D

Lab ID: 240-59503-G-4-C MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Aroclor-1016 | 441 | 148 | 33 | 17 | 56 | 17-141 | |
| Aroclor-1260 | 441 | 118 | 27 | 29 | 55 | 19-137 | p |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM IV
PCBS METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: MB 240-212544/7-A

Matrix: Solid  Date Extracted: 12/30/2015 06:54

Lab File ID:(1)  Lab File ID:(2) P1200019.D

Date Analyzed:(1)  Date Analyzed:(2) 01/04/2016 12:43

Instrument ID:(1)  Instrument ID:(2) A2HP12

GC Column:(1)  ID:  GC Column:(2) CLP-1  ID: 0.53(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| F-VB-UST-P-7 | 240-59482-2 | | 01/04/2016 10:19 |
| F-VB-UST-P-8 | 240-59482-3 | | 01/04/2016 10:37 |
| | 240-59503-G-4-B MS | | 01/04/2016 12:07 |
| | 240-59503-G-4-C MSD | | 01/04/2016 12:25 |
| | LCS 240-212544/8-A | | 01/04/2016 13:01 |

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Sample No.: STD05 240-207007/31  Date Analyzed: 11/17/2015 04:23

Instrument ID: A2HP12  GC Column: CLP-1  ID: 0.53(mm)

Lab File ID (Standard): P1200066.D  Heated Purge: (Y/N) N

Calibration ID: 31555

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 188112954 | 2.68 | | | | |
| UPPER LIMIT | | 376225908 | 3.18 | | | | |
| LOWER LIMIT | | 94056477 | 2.18 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-207007/39 | | 184850376 | 2.69 | | | | |
| CCV 240-212845/3 CCVIS | | 201292307 | 2.69 | | | | |

BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Sample No.: CCV 240-212845/3 Date Analyzed: 01/04/2016 07:42

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm)

Lab File ID (Standard): P1200003.D Heated Purge: (Y/N) N

Calibration ID: 31555

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 201292307 | 2.69 | | | | |
| UPPER LIMIT | | 402584614 | 3.19 | | | | |
| LOWER LIMIT | | 100646154 | 2.19 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-212845/4 | | 193760566 | 2.69 | | | | |
| CCV 240-212845/5 | | 196884227 | 2.69 | | | | |
| CCV 240-212845/6 | | 201868653 | 2.69 | | | | |
| CCV 240-212845/7 | | 187655215 | 2.69 | | | | |
| 240-59482-2 | F-VB-UST-P-7 | 208904735 | 2.70 | | | | |
| 240-59482-3 | F-VB-UST-P-8 | 184367355 | 2.69 | | | | |
| 240-59503-G-4-B MS | | 229769056 | 2.69 | | | | |
| 240-59503-G-4-C MSD | | 314864953 | 2.69 | | | | |
| MB 240-212544/7-A | | 199821577 | 2.69 | | | | |
| LCS 240-212544/8-A | | 199206956 | 2.70 | | | | |

BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM X
IDENTIFICATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Client Sample ID: Lab Sample ID: 240-59503-G-4-B MS

Instrument ID (1): Instrument ID (2): A2HP12

Date Analyzed (1): Date Analyzed (2): 01/04/2016 12:07

GC Column (1): ID: GC Column (2): CLP-1 ID: 0.53(mm)

| ANALYTE | COL | PEAK | RT | RT WINDOW | | CONCENTRATION | | |
|---|---|---|---|---|---|---|---|---|
| | | | | FROM | TO | PEAK | MEAN | |
| Aroclor-1016 | 1 | 1 | 4.61 | 4.58 | 4.64 | 233 | 175 | |
| | | 2 | 5.11 | 5.08 | 5.14 | 200 | | |
| | | 3 | 5.72 | 5.71 | 5.77 | 94.1 | | |
| | | 4 | 5.91 | 5.88 | 5.94 | 165 | | |
| | | 5 | 6.45 | 6.42 | 6.48 | 183 | | |
| Aroclor-1260 | 1 | 1 | 7.95 | 7.92 | 7.98 | 156 | 143 | |
| | | 2 | 8.28 | 8.26 | 8.32 | 149 | | |
| | | 3 | 8.61 | 8.58 | 8.64 | 137 | | |
| | | 4 | 9.40 | 9.37 | 9.43 | 138 | | |
| | | 5 | 9.71 | 9.69 | 9.75 | 136 | | |

FORM X
IDENTIFICATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Client Sample ID: Lab Sample ID: 240-59503-G-4-C MSD

Instrument ID (1): Instrument ID (2): A2HP12

Date Analyzed (1): Date Analyzed (2): 01/04/2016 12:25

GC Column (1): ID: GC Column (2): CLP-1 ID: 0.53(mm)

| ANALYTE | COL | PEAK | RT | RT WINDOW | | CONCENTRATION | | |
|---|---|---|---|---|---|---|---|---|
| | | | | FROM | TO | PEAK | MEAN | |
| Aroclor-1016 | 1 | 1 | 4.61 | 4.58 | 4.64 | 241 | 148 | |
| | | 2 | 5.11 | 5.08 | 5.14 | 135 | | |
| | | 3 | 5.73 | 5.71 | 5.77 | 113 | | |
| | | 4 | 5.91 | 5.88 | 5.94 | 119 | | |
| | | 5 | 6.44 | 6.42 | 6.48 | 130 | | |
| Aroclor-1260 | 1 | 1 | 7.95 | 7.92 | 7.98 | 127 | 118 | |
| | | 2 | 8.28 | 8.26 | 8.32 | 120 | | |
| | | 3 | 8.61 | 8.58 | 8.64 | 113 | | |
| | | 4 | 9.39 | 9.37 | 9.43 | 119 | | |
| | | 5 | 9.71 | 9.69 | 9.75 | 113 | | |

FORM X
IDENTIFICATION SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Client Sample ID: Lab Sample ID: LCS 240-212544/8-A

Instrument ID (1): Instrument ID (2): A2HP12

Date Analyzed (1): Date Analyzed (2): 01/04/2016 13:01

GC Column (1): ID: GC Column (2): CLP-1 ID: 0.53(mm)

| ANALYTE | COL | PEAK | RT | RT WINDOW | | CONCENTRATION | | |
|---|---|---|---|---|---|---|---|---|
| | | | | FROM | TO | PEAK | MEAN | |
| Aroclor-1016 | 1 | 1 | 4.61 | 4.58 | 4.64 | 326 | 314 | |
| | | 2 | 5.12 | 5.08 | 5.14 | 326 | | |
| | | 3 | 5.74 | 5.71 | 5.77 | 269 | | |
| | | 4 | 5.92 | 5.88 | 5.94 | 332 | | |
| | | 5 | 6.45 | 6.42 | 6.48 | 319 | | |
| Aroclor-1260 | 1 | 1 | 7.95 | 7.92 | 7.98 | 332 | 337 | |
| | | 2 | 8.29 | 8.26 | 8.32 | 336 | | |
| | | 3 | 8.61 | 8.58 | 8.64 | 315 | | |
| | | 4 | 9.40 | 9.37 | 9.43 | 370 | | |
| | | 5 | 9.72 | 9.69 | 9.75 | 333 | | |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-7 | Lab Sample ID: 240-59482-2

Matrix: Solid | Lab File ID: P1200011.D

Analysis Method: 8082A | Date Collected: 12/28/2015 10:00

Extraction Method: 3540C | Date Extracted: 12/30/2015 06:54

Sample wt/vol: 29.83(g) | Date Analyzed: 01/04/2016 10:19

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 18.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212845 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 15 | U | 41 | 15 |
| 11104-28-2 | Aroclor-1221 | 20 | U | 41 | 20 |
| 11141-16-5 | Aroclor-1232 | 25 | U | 41 | 25 |
| 53469-21-9 | Aroclor-1242 | 14 | U | 41 | 14 |
| 12672-29-6 | Aroclor-1248 | 9.9 | U | 41 | 9.9 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 41 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 41 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 84 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 72 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: F-VB-UST-P-8 | Lab Sample ID: 240-59482-3

Matrix: Solid | Lab File ID: P1200012.D

Analysis Method: 8082A | Date Collected: 12/28/2015 10:30

Extraction Method: 3540C | Date Extracted: 12/30/2015 06:54

Sample wt/vol: 29.77(g) | Date Analyzed: 01/04/2016 10:37

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 16.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212845 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 40 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.6 | U | 40 | 9.6 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 57 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 92 | | 10-155 |

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 20:14 Calibration End Date: 11/16/2015 21:45 Calibration ID: 31527

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/4 | P1200039.D |
| Level 2 | STD01 240-207007/5 | P1200040.D |
| Level 3 | STD02 240-207007/6 | P1200041.D |
| Level 4 | STD05 240-207007/7 | P1200042.D |
| Level 5 | STD1 240-207007/8 | P1200043.D |
| Level 6 | STD2 240-207007/9 | P1200044.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1232 Peak 1 | 0.0146 0.0120 | 0.0146 | 0.0137 | 0.0137 | 0.0123 | Lin1 | 0.0002 | 0.0122 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 2 | 0.0104 0.0088 | 0.0106 | 0.0101 | 0.0100 | 0.0090 | Lin1 | 0.0001 | 0.0089 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 3 | 0.0214 0.0206 | 0.0218 | 0.0204 | 0.0219 | 0.0205 | Lin1 | 0.0001 | 0.0207 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1232 Peak 4 | 0.0088 0.0084 | 0.0090 | 0.0081 | 0.0093 | 0.0085 | Lin1 | 0 | 0.0085 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 5 | 0.0088 0.0076 | 0.0092 | 0.0087 | 0.0086 | 0.0078 | Lin1 | 0.0001 | 0.0078 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1262 Peak 1 | 0.0288 0.0238 | 0.0249 | 0.0236 | 0.0241 | 0.0228 | Lin1 | 0.0002 | 0.0233 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1262 Peak 2 | 0.0420 0.0453 | 0.0443 | 0.0426 | 0.0462 | 0.0441 | Lin1 | 0 | 0.0452 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 3 | 0.0380 0.0415 | 0.0404 | 0.0397 | 0.0426 | 0.0405 | Lin1 | 0 | 0.0415 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 4 | 0.0853 0.0976 | 0.0928 | 0.0961 | 0.0994 | 0.0962 | Lin1 | -0.001 | 0.0979 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 5 | 0.0364 0.0385 | 0.0386 | 0.0374 | 0.0395 | 0.0372 | Lin1 | 0 | 0.0383 | | | | | | | 0.9990 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 20:14 Calibration End Date: 11/16/2015 21:45 Calibration ID: 31527

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/4 | P1200039.D |
| Level 2 | STD01 240-207007/5 | P1200040.D |
| Level 3 | STD02 240-207007/6 | P1200041.D |
| Level 4 | STD05 240-207007/7 | P1200042.D |
| Level 5 | STD1 240-207007/8 | P1200043.D |
| Level 6 | STD2 240-207007/9 | P1200044.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1232 Peak 1 | BNB | Lin1 | 2864462<br>94613461 | 5743911 | 10579052 | 25148552 | 48682811 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 2 | BNB | Lin1 | 2042318<br>69008204 | 4176797 | 7799621 | 18338381 | 35520928 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 3 | BNB | Lin1 | 4204092<br>162234139 | 8567857 | 15765426 | 40096819 | 80981509 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 4 | BNB | Lin1 | 1732949<br>65818756 | 3551392 | 6289163 | 17045553 | 33804666 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1232 Peak 5 | BNB | Lin1 | 1725581<br>59962732 | 3622594 | 6723266 | 15858414 | 31021460 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 1 | BNB | Lin1 | 5662298<br>186973877 | 9773527 | 18252707 | 44264730 | 89983608 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 2 | BNB | Lin1 | 8249286<br>356026718 | 17431705 | 32928458 | 84853483 | 174547806 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 3 | BNB | Lin1 | 7469803<br>326504754 | 15875631 | 30684939 | 78118840 | 160322518 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 4 | BNB | Lin1 | 16781995<br>766924713 | 36487355 | 74250353 | 182460886 | 380317708 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1262 Peak 5 | BNB | Lin1 | 7158889<br>302670598 | 15182112 | 28944443 | 72530686 | 147149653 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 22:03 Calibration End Date: 11/16/2015 23:33 Calibration ID: 31533

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/10 | P1200045.D |
| Level 2 | STD01 240-207007/11 | P1200046.D |
| Level 3 | STD02 240-207007/12 | P1200047.D |
| Level 4 | STD05 240-207007/13 | P1200048.D |
| Level 5 | STD1 240-207007/14 | P1200049.D |
| Level 6 | STD2 240-207007/15 | P1200050.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1242 Peak 1 | 0.0115 0.0082 | 0.0100 | 0.0094 | 0.0091 | 0.0087 | Lin1 | 0.0002 | 0.0083 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 2 | 0.0201 0.0153 | 0.0180 | 0.0172 | 0.0169 | 0.0161 | Lin1 | 0.0003 | 0.0155 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 3 | 0.0408 0.0378 | 0.0334 | 0.0373 | 0.0380 | 0.0380 | Lin1 | 0 | 0.0379 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 4 | 0.0176 0.0154 | 0.0131 | 0.0162 | 0.0160 | 0.0157 | Lin1 | 0 | 0.0155 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1242 Peak 5 | 0.0196 0.0154 | 0.0173 | 0.0163 | 0.0163 | 0.0157 | Lin1 | 0.0002 | 0.0155 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 1 | 0.1160 0.1140 | 0.1051 | 0.1050 | 0.1121 | 0.1102 | Lin1 | 0 | 0.1126 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1268 Peak 2 | 0.1129 0.1102 | 0.1035 | 0.1031 | 0.1082 | 0.1086 | Lin1 | 0 | 0.1094 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 3 | 0.0941 0.0941 | 0.0865 | 0.0873 | 0.0935 | 0.0918 | Lin1 | 0 | 0.0934 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 4 | 0.0427 0.0386 | 0.0375 | 0.0368 | 0.0384 | 0.0379 | Lin1 | 0.0001 | 0.0382 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 5 | 0.3010 0.3253 | 0.2825 | 0.2850 | 0.3104 | 0.3105 | Lin1 | -0.003 | 0.3204 | | | | | | | 0.9990 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 22:03 Calibration End Date: 11/16/2015 23:33 Calibration ID: 31533

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/10 | P1200045.D |
| Level 2 | STD01 240-207007/11 | P1200046.D |
| Level 3 | STD02 240-207007/12 | P1200047.D |
| Level 4 | STD05 240-207007/13 | P1200048.D |
| Level 5 | STD1 240-207007/14 | P1200049.D |
| Level 6 | STD2 240-207007/15 | P1200050.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1242 Peak 1 | BNB | Lin1 | 2149767<br>64172683 | 3972268 | 7542438 | 18026385 | 33712614 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 2 | BNB | Lin1 | 3777832<br>120095985 | 7167048 | 13831308 | 33547847 | 62739379 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 3 | BNB | Lin1 | 7649911<br>296721464 | 13315192 | 29900733 | 75211737 | 147999390 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 4 | BNB | Lin1 | 3299090<br>120600215 | 5223145 | 12995644 | 31796203 | 60982553 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1242 Peak 5 | BNB | Lin1 | 3672196<br>121118126 | 6903841 | 13110974 | 32386776 | 61129587 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 1 | BNB | Lin1 | 21759307<br>894333238 | 41863019 | 84230109 | 222130676 | 428635453 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 2 | BNB | Lin1 | 21178393<br>864175846 | 41209432 | 82708339 | 214470756 | 422311987 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 3 | BNB | Lin1 | 17646266<br>738380844 | 34461725 | 70030810 | 185306823 | 357127789 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 4 | BNB | Lin1 | 8003779<br>302725873 | 14951207 | 29550936 | 76139152 | 147585310 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1268 Peak 5 | BNB | Lin1 | 56470647<br>2551765384 | 112539547 | 228706125 | 615081152 | 1208047688 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 23:51 Calibration End Date: 11/17/2015 01:22 Calibration ID: 31539

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/16 | P1200051.D |
| Level 2 | STD01 240-207007/17 | P1200052.D |
| Level 3 | STD02 240-207007/18 | P1200053.D |
| Level 4 | STD05 240-207007/19 | P1200054.D |
| Level 5 | STD1 240-207007/20 | P1200055.D |
| Level 6 | STD2 240-207007/21 | P1200056.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1248 Peak 1 | 0.0111 0.0082 | 0.0099 | 0.0099 | 0.0090 | 0.0084 | Lin1 | 0.0002 | 0.0082 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1248 Peak 2 | 0.0277 0.0242 | 0.0256 | 0.0251 | 0.0243 | 0.0239 | Lin1 | 0.0002 | 0.0240 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 3 | 0.0240 0.0225 | 0.0232 | 0.0233 | 0.0226 | 0.0219 | Lin1 | 0.0001 | 0.0223 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 4 | 0.0199 0.0193 | 0.0194 | 0.0192 | 0.0192 | 0.0190 | Lin1 | 0 | 0.0192 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 5 | 0.0138 0.0128 | 0.0122 | 0.0128 | 0.0125 | 0.0116 | Lin1 | 0 | 0.0124 | | | | | | | 0.9980 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/16/2015 23:51 Calibration End Date: 11/17/2015 01:22 Calibration ID: 31539

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/16 | P1200051.D |
| Level 2 | STD01 240-207007/17 | P1200052.D |
| Level 3 | STD02 240-207007/18 | P1200053.D |
| Level 4 | STD05 240-207007/19 | P1200054.D |
| Level 5 | STD1 240-207007/20 | P1200055.D |
| Level 6 | STD2 240-207007/21 | P1200056.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1248 Peak 1 | BNB | Lin1 | 2214671<br>65344382 | 3893329 | 7710225 | 17465613 | 33070601 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 2 | BNB | Lin1 | 5540374<br>193008938 | 10066768 | 19570159 | 46979815 | 94056087 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 3 | BNB | Lin1 | 4803789<br>179213673 | 9149673 | 18179130 | 43818711 | 86223823 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 4 | BNB | Lin1 | 3975067<br>153520953 | 7642679 | 14994793 | 37207246 | 74584605 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1248 Peak 5 | BNB | Lin1 | 2764694<br>102029173 | 4797959 | 10009264 | 24248210 | 45732697 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/17/2015 01:40 Calibration End Date: 11/17/2015 03:10 Calibration ID: 31549

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/22 | P1200057.D |
| Level 2 | STD01 240-207007/23 | P1200058.D |
| Level 3 | STD02 240-207007/24 | P1200059.D |
| Level 4 | STD05 240-207007/25 | P1200060.D |
| Level 5 | STD1 240-207007/26 | P1200061.D |
| Level 6 | STD2 240-207007/27 | P1200062.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1221 Peak 1 | 0.0083 0.0057 | 0.0074 | 0.0067 | 0.0065 | 0.0061 | Lin1 | 0.0002 | 0.0058 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1221 Peak 2 | 0.0055 0.0037 | 0.0049 | 0.0045 | 0.0042 | 0.0039 | Lin1 | 0.0001 | 0.0037 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1221 Peak 3 | 0.0208 0.0145 | 0.0183 | 0.0167 | 0.0164 | 0.0154 | Lin1 | 0.0004 | 0.0147 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1254 Peak 1 | 0.0259 0.0202 | 0.0228 | 0.0210 | 0.0212 | 0.0203 | Lin1 | 0.0003 | 0.0201 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 2 | 0.0354 0.0290 | 0.0313 | 0.0287 | 0.0298 | 0.0292 | Lin1 | 0.0003 | 0.0289 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 3 | 0.0466 0.0419 | 0.0442 | 0.0413 | 0.0427 | 0.0391 | Lin1 | 0.0003 | 0.0410 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1254 Peak 4 | 0.0442 0.0360 | 0.0392 | 0.0345 | 0.0358 | 0.0355 | Lin1 | 0.0003 | 0.0354 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1254 Peak 5 | 0.0544 0.0459 | 0.0486 | 0.0448 | 0.0465 | 0.0453 | Lin1 | 0.0003 | 0.0454 | | | | | | | 1.0000 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/17/2015 01:40 Calibration End Date: 11/17/2015 03:10 Calibration ID: 31549

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/22 | P1200057.D |
| Level 2 | STD01 240-207007/23 | P1200058.D |
| Level 3 | STD02 240-207007/24 | P1200059.D |
| Level 4 | STD05 240-207007/25 | P1200060.D |
| Level 5 | STD1 240-207007/26 | P1200061.D |
| Level 6 | STD2 240-207007/27 | P1200062.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1221 Peak 1 | BNB | Lin1 | 1617753<br>43159048 | 2834146 | 5288305 | 12630620 | 23444359 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1221 Peak 2 | BNB | Lin1 | 1073929<br>27877416 | 1853226 | 3515548 | 8239497 | 15092745 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1221 Peak 3 | BNB | Lin1 | 4061658<br>109234720 | 6958986 | 13214576 | 31899214 | 59455965 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 1 | BNB | Lin1 | 5070982<br>152082598 | 8694593 | 16601983 | 41262994 | 78538732 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 2 | BNB | Lin1 | 6915782<br>218988362 | 11916252 | 22685227 | 57976400 | 113180990 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 3 | BNB | Lin1 | 9122282<br>316127715 | 16841606 | 32606706 | 82955431 | 151535520 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 4 | BNB | Lin1 | 8646084<br>271574308 | 14939001 | 27264214 | 69542078 | 137323388 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1254 Peak 5 | BNB | Lin1 | 10643906<br>346633557 | 18530884 | 35338412 | 90448562 | 175425464 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/17/2015 03:29 Calibration End Date: 11/17/2015 04:59 Calibration ID: 31555

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/28 | P1200063.D |
| Level 2 | STD01 240-207007/29 | P1200064.D |
| Level 3 | STD02 240-207007/30 | P1200065.D |
| Level 4 | STD05 240-207007/31 | P1200066.D |
| Level 5 | STD1 240-207007/32 | P1200067.D |
| Level 6 | STD2 240-207007/33 | P1200068.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1016 Peak 1 | 0.0130 0.0104 | 0.0132 | 0.0125 | 0.0117 | 0.0109 | Lin1 | 0.0002 | 0.0106 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1016 Peak 2 | 0.0226 0.0190 | 0.0238 | 0.0222 | 0.0210 | 0.0197 | Lin1 | 0.0003 | 0.0193 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1016 Peak 3 | 0.0471 0.0469 | 0.0503 | 0.0465 | 0.0479 | 0.0469 | Lin1 | 0.0001 | 0.0469 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 4 | 0.0212 0.0189 | 0.0222 | 0.0202 | 0.0199 | 0.0193 | Lin1 | 0.0002 | 0.0190 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 5 | 0.0220 0.0187 | 0.0228 | 0.0210 | 0.0199 | 0.0191 | Lin1 | 0.0003 | 0.0188 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1260 Peak 1 | 0.0330 0.0301 | 0.0338 | 0.0324 | 0.0309 | 0.0302 | Lin1 | 0.0003 | 0.0300 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 2 | 0.0622 0.0585 | 0.0618 | 0.0622 | 0.0606 | 0.0586 | Lin1 | 0.0003 | 0.0586 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 3 | 0.0650 0.0628 | 0.0634 | 0.0652 | 0.0643 | 0.0626 | Lin1 | 0.0002 | 0.0629 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 4 | 0.0661 0.0714 | 0.0711 | 0.0771 | 0.0773 | 0.0731 | Lin1 | 0 | 0.0729 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1260 Peak 5 | 0.0374 0.0384 | 0.0399 | 0.0382 | 0.0386 | 0.0385 | Lin1 | 0 | 0.0384 | | | | | | | 1.0000 | | 0.9900 |
| Tetrachloro-m-xylene | 0.7048 0.6608 | 0.7003 | 0.7144 | 0.6986 | 0.6822 | Lin1 | 0.0002 | 0.6699 | | | | | | | 0.9990 | | 0.9900 |
| DCB Decachlorobiphenyl | 0.7699 0.6896 | 0.7629 | 0.7446 | 0.7316 | 0.6953 | Lin1 | 0.0003 | 0.6927 | | | | | | | 1.0000 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Canton Job No.: 240-59482-1 Analy Batch No.: 207007

SDG No.:

Instrument ID: A2HP12 GC Column: CLP-1 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/17/2015 03:29 Calibration End Date: 11/17/2015 04:59 Calibration ID: 31555

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-207007/28 | P1200063.D |
| Level 2 | STD01 240-207007/29 | P1200064.D |
| Level 3 | STD02 240-207007/30 | P1200065.D |
| Level 4 | STD05 240-207007/31 | P1200066.D |
| Level 5 | STD1 240-207007/32 | P1200067.D |
| Level 6 | STD2 240-207007/33 | P1200068.D |

| ANALYTE | IS REF | CURVE TYPE | RESPONSE | | | | | CONCENTRATION (NG/UL) | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 |
| PCB-1016 Peak 1 | BNB | Lin1 | 2707871<br>80765612 | 4820722 | 9085846 | 22006625 | 42447741 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 2 | BNB | Lin1 | 4700873<br>147609677 | 8734684 | 16183493 | 39469767 | 76715703 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 3 | BNB | Lin1 | 9792259<br>364184674 | 18438138 | 33886380 | 90045825 | 182657834 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 4 | BNB | Lin1 | 4396897<br>147122332 | 8128551 | 14710049 | 37512861 | 75171974 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1016 Peak 5 | BNB | Lin1 | 4563462<br>145010499 | 8352402 | 15341259 | 37426880 | 74524891 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 1 | BNB | Lin1 | 6849336<br>233433821 | 12383500 | 23597918 | 58080604 | 117631104 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 2 | BNB | Lin1 | 12922388<br>454390240 | 22631342 | 45372536 | 113933482 | 228581266 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 3 | BNB | Lin1 | 13513426<br>488067042 | 23245320 | 47522931 | 120987145 | 244049421 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 4 | BNB | Lin1 | 13737456<br>554191846 | 26050869 | 56208655 | 145378951 | 285103725 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| PCB-1260 Peak 5 | BNB | Lin1 | 7764692<br>298012472 | 14617316 | 27852633 | 72613053 | 149980149 | 0.0500<br>2.00 | 0.100 | 0.200 | 0.500 | 1.00 |
| Tetrachloro-m-xylene | BNB | Lin1 | 7322625<br>256611202 | 12827811 | 26051804 | 65705303 | 132973208 | 0.00250<br>0.100 | 0.00500 | 0.0100 | 0.0250 | 0.0500 |
| DCB Decachlorobiphenyl | BNB | Lin1 | 7998062<br>267771950 | 13974803 | 27151934 | 68816182 | 135510483 | 0.00250<br>0.100 | 0.00500 | 0.0100 | 0.0250 | 0.0500 |

Curve Type Legend:
Lin1 = Linear 1/conc ISTD

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-207007/39  Calibration Date: 11/17/2015 06:48

Instrument ID: A2HP12  Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/17/2015 04:59

Lab File ID: P1200074.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0109 | | 1.01 | 1.00 | 1.2 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0209 | | 1.07 | 1.00 | 6.9 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0504 | | 1.07 | 1.00 | 7.2 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0205 | | 1.07 | 1.00 | 6.9 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0196 | | 1.03 | 1.00 | 3.0 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0302 | | 0.995 | 1.00 | -0.5 | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0570 | | 0.967 | 1.00 | -3.3 | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0535 | | 0.848 | 1.00 | -15.2 | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0824 | | 1.13 | 1.00 | 13.1 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0405 | | 1.05 | 1.00 | 5.3 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6836 | | 0.0507 | 0.0500 | 1.5 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.6921 | | 0.0495 | 0.0500 | -1.0 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: ICV 240-207007/39 Calibration Date: 11/17/2015 06:48

Instrument ID: A2HP12 Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 04:59

Lab File ID: P1200074.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1016 Peak 1 | 4.61 | 4.58 | 4.64 |
| PCB-1016 Peak 2 | 5.11 | 5.08 | 5.14 |
| PCB-1016 Peak 3 | 5.74 | 5.71 | 5.77 |
| PCB-1016 Peak 4 | 5.91 | 5.88 | 5.94 |
| PCB-1016 Peak 5 | 6.45 | 6.42 | 6.48 |
| PCB-1260 Peak 1 | 7.95 | 7.92 | 7.98 |
| PCB-1260 Peak 2 | 8.29 | 8.26 | 8.32 |
| PCB-1260 Peak 3 | 8.61 | 8.58 | 8.64 |
| PCB-1260 Peak 4 | 9.40 | 9.37 | 9.43 |
| PCB-1260 Peak 5 | 9.72 | 9.69 | 9.75 |
| Tetrachloro-m-xylene | 4.05 | 4.02 | 4.08 |
| DCB Decachlorobiphenyl | 10.93 | 10.90 | 10.96 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/3 Calibration Date: 01/04/2016 07:42

Instrument ID: A2HP12 Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 04:59

Lab File ID: P1200003.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0125 | | 0.569 | 0.500 | 13.8 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0222 | | 0.558 | 0.500 | 11.5 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0487 | | 0.517 | 0.500 | 3.3 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0209 | | 0.539 | 0.500 | 7.8 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0207 | | 0.535 | 0.500 | 7.1 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0331 | | 0.541 | 0.500 | 8.3 | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0635 | | 0.536 | 0.500 | 7.1 | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0671 | | 0.531 | 0.500 | 6.2 | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0813 | | 0.558 | 0.500 | 11.6 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0395 | | 0.513 | 0.500 | 2.7 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.7403 | | 0.0273 | 0.0250 | 9.4 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.7093 | | 0.0252 | 0.0250 | 0.6 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/3 Calibration Date: 01/04/2016 07:42

Instrument ID: A2HP12 Calib Start Date: 11/17/2015 03:29

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 04:59

Lab File ID: P1200003.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1016 Peak 1 | 4.61 | 4.58 | 4.64 |
| PCB-1016 Peak 2 | 5.11 | 5.08 | 5.14 |
| PCB-1016 Peak 3 | 5.74 | 5.71 | 5.77 |
| PCB-1016 Peak 4 | 5.91 | 5.88 | 5.94 |
| PCB-1016 Peak 5 | 6.45 | 6.42 | 6.48 |
| PCB-1260 Peak 1 | 7.95 | 7.92 | 7.98 |
| PCB-1260 Peak 2 | 8.29 | 8.26 | 8.32 |
| PCB-1260 Peak 3 | 8.61 | 8.58 | 8.64 |
| PCB-1260 Peak 4 | 9.40 | 9.37 | 9.43 |
| PCB-1260 Peak 5 | 9.72 | 9.69 | 9.75 |
| Tetrachloro-m-xylene | 4.05 | 4.02 | 4.08 |
| DCB Decachlorobiphenyl | 10.92 | 10.89 | 10.95 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/4  Calibration Date: 01/04/2016 08:00

Instrument ID: A2HP12  Calib Start Date: 11/16/2015 20:14

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/16/2015 21:45

Lab File ID: P1200004.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1232 Peak 1 | Lin1 | | 0.0149 | | 0.596 | 0.500 | 19.1 | 20.0 |
| PCB-1232 Peak 2 | Lin1 | | 0.0107 | | 0.585 | 0.500 | 17.0 | 20.0 |
| PCB-1232 Peak 3 | Lin1 | | 0.0217 | | 0.521 | 0.500 | 4.2 | 20.0 |
| PCB-1232 Peak 4 | Lin1 | | 0.0093 | | 0.542 | 0.500 | 8.4 | 20.0 |
| PCB-1232 Peak 5 | Lin1 | | 0.0092 | | 0.575 | 0.500 | 15.1 | 20.0 |
| PCB-1262 Peak 1 | Lin1 | | 0.0258 | | 0.545 | 0.500 | 9.0 | 20.0 |
| PCB-1262 Peak 2 | Lin1 | | 0.0457 | | 0.510 | 0.500 | 2.0 | 20.0 |
| PCB-1262 Peak 3 | Lin1 | | 0.0438 | | 0.532 | 0.500 | 6.4 | 20.0 |
| PCB-1262 Peak 4 | Lin1 | | 0.1069 | | 0.551 | 0.500 | 10.3 | 20.0 |
| PCB-1262 Peak 5 | Lin1 | | 0.0421 | | 0.551 | 0.500 | 10.1 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/4 Calibration Date: 01/04/2016 08:00

Instrument ID: A2HP12 Calib Start Date: 11/16/2015 20:14

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/16/2015 21:45

Lab File ID: P1200004.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1232 Peak 1 | 4.61 | 4.58 | 4.64 |
| PCB-1232 Peak 2 | 5.11 | 5.08 | 5.14 |
| PCB-1232 Peak 3 | 5.74 | 5.71 | 5.77 |
| PCB-1232 Peak 4 | 5.91 | 5.88 | 5.94 |
| PCB-1232 Peak 5 | 6.45 | 6.42 | 6.48 |
| PCB-1262 Peak 1 | 7.95 | 7.92 | 7.98 |
| PCB-1262 Peak 2 | 8.75 | 8.72 | 8.78 |
| PCB-1262 Peak 3 | 9.03 | 9.00 | 9.06 |
| PCB-1262 Peak 4 | 9.40 | 9.37 | 9.43 |
| PCB-1262 Peak 5 | 9.77 | 9.74 | 9.80 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/5  Calibration Date: 01/04/2016 08:19

Instrument ID: A2HP12  Calib Start Date: 11/16/2015 22:03

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/16/2015 23:33

Lab File ID: P1200005.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1242 Peak 1 | Lin1 | | 0.0093 | | 0.541 | 0.500 | 8.2 | 20.0 |
| PCB-1242 Peak 2 | Lin1 | | 0.0173 | | 0.540 | 0.500 | 8.1 | 20.0 |
| PCB-1242 Peak 3 | Lin1 | | 0.0374 | | 0.495 | 0.500 | -0.9 | 20.0 |
| PCB-1242 Peak 4 | Lin1 | | 0.0161 | | 0.518 | 0.500 | 3.6 | 20.0 |
| PCB-1242 Peak 5 | Lin1 | | 0.0166 | | 0.522 | 0.500 | 4.4 | 20.0 |
| PCB-1268 Peak 1 | Lin1 | | 0.1106 | | 0.494 | 0.500 | -1.2 | 20.0 |
| PCB-1268 Peak 2 | Lin1 | | 0.1122 | | 0.515 | 0.500 | 3.0 | 20.0 |
| PCB-1268 Peak 3 | Lin1 | | 0.0943 | | 0.508 | 0.500 | 1.7 | 20.0 |
| PCB-1268 Peak 4 | Lin1 | | 0.0381 | | 0.496 | 0.500 | -0.8 | 20.0 |
| PCB-1268 Peak 5 | Lin1 | | 0.3053 | | 0.485 | 0.500 | -3.0 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/5 Calibration Date: 01/04/2016 08:19

Instrument ID: A2HP12 Calib Start Date: 11/16/2015 22:03

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/16/2015 23:33

Lab File ID: P1200005.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1242 Peak 1 | 4.61 | 4.58 | 4.64 |
| PCB-1242 Peak 2 | 5.11 | 5.08 | 5.14 |
| PCB-1242 Peak 3 | 5.74 | 5.71 | 5.77 |
| PCB-1242 Peak 4 | 5.92 | 5.89 | 5.95 |
| PCB-1242 Peak 5 | 6.45 | 6.42 | 6.48 |
| PCB-1268 Peak 1 | 9.77 | 9.74 | 9.80 |
| PCB-1268 Peak 2 | 9.81 | 9.78 | 9.84 |
| PCB-1268 Peak 3 | 10.03 | 10.00 | 10.06 |
| PCB-1268 Peak 4 | 10.44 | 10.41 | 10.47 |
| PCB-1268 Peak 5 | 10.74 | 10.71 | 10.77 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/6  Calibration Date: 01/04/2016 08:37

Instrument ID: A2HP12  Calib Start Date: 11/16/2015 23:51

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/17/2015 01:22

Lab File ID: P1200006.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1248 Peak 1 | Lin1 | | 0.0100 | | 0.585 | 0.500 | 17.0 | 20.0 |
| PCB-1248 Peak 2 | Lin1 | | 0.0245 | | 0.503 | 0.500 | 0.5 | 20.0 |
| PCB-1248 Peak 3 | Lin1 | | 0.0222 | | 0.493 | 0.500 | -1.5 | 20.0 |
| PCB-1248 Peak 4 | Lin1 | | 0.0194 | | 0.504 | 0.500 | 0.9 | 20.0 |
| PCB-1248 Peak 5 | Lin1 | | 0.0123 | | 0.492 | 0.500 | -1.6 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/6 Calibration Date: 01/04/2016 08:37

Instrument ID: A2HP12 Calib Start Date: 11/16/2015 23:51

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 01:22

Lab File ID: P1200006.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1248 Peak 1 | 5.11 | 5.08 | 5.14 |
| PCB-1248 Peak 2 | 5.74 | 5.71 | 5.77 |
| PCB-1248 Peak 3 | 7.01 | 6.98 | 7.04 |
| PCB-1248 Peak 4 | 7.25 | 7.22 | 7.28 |
| PCB-1248 Peak 5 | 7.73 | 7.70 | 7.76 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/7  Calibration Date: 01/04/2016 08:55

Instrument ID: A2HP12  Calib Start Date: 11/17/2015 01:40

GC Column: CLP-1  ID: 0.53(mm)  Calib End Date: 11/17/2015 03:10

Lab File ID: P1200007.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1221 Peak 1 | Lin1 | | 0.0069 | | 0.565 | 0.500 | 13.0 | 20.0 |
| PCB-1221 Peak 2 | Lin1 | | 0.0043 | | 0.547 | 0.500 | 9.4 | 20.0 |
| PCB-1221 Peak 3 | Lin1 | | 0.0173 | | 0.564 | 0.500 | 12.9 | 20.0 |
| PCB-1254 Peak 1 | Lin1 | | 0.0220 | | 0.534 | 0.500 | 6.8 | 20.0 |
| PCB-1254 Peak 2 | Lin1 | | 0.0302 | | 0.513 | 0.500 | 2.6 | 20.0 |
| PCB-1254 Peak 3 | Lin1 | | 0.0394 | | 0.474 | 0.500 | -5.1 | 20.0 |
| PCB-1254 Peak 4 | Lin1 | | 0.0335 | | 0.464 | 0.500 | -7.2 | 20.0 |
| PCB-1254 Peak 5 | Lin1 | | 0.0462 | | 0.501 | 0.500 | 0.2 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Lab Sample ID: CCV 240-212845/7 Calibration Date: 01/04/2016 08:55

Instrument ID: A2HP12 Calib Start Date: 11/17/2015 01:40

GC Column: CLP-1 ID: 0.53(mm) Calib End Date: 11/17/2015 03:10

Lab File ID: P1200007.D

| Analyte | RT | RT WINDOW | |
|---|---|---|---|
| | | FROM | TO |
| PCB-1221 Peak 1 | 4.34 | 4.31 | 4.37 |
| PCB-1221 Peak 2 | 4.55 | 4.52 | 4.58 |
| PCB-1221 Peak 3 | 4.61 | 4.58 | 4.64 |
| PCB-1254 Peak 1 | 6.93 | 6.90 | 6.96 |
| PCB-1254 Peak 2 | 7.26 | 7.23 | 7.29 |
| PCB-1254 Peak 3 | 7.73 | 7.70 | 7.76 |
| PCB-1254 Peak 4 | 8.07 | 8.04 | 8.10 |
| PCB-1254 Peak 5 | 8.61 | 8.58 | 8.64 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-212544/7-A

Matrix: Solid | Lab File ID: P1200019.D

Analysis Method: 8082A | Date Collected:

Extraction Method: 3540C | Date Extracted: 12/30/2015 06:54

Sample wt/vol: 30(g) | Date Analyzed: 01/04/2016 12:43

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212845 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 12 | U | 33 | 12 |
| 11104-28-2 | Aroclor-1221 | 16 | U | 33 | 16 |
| 11141-16-5 | Aroclor-1232 | 20 | U | 33 | 20 |
| 53469-21-9 | Aroclor-1242 | 11 | U | 33 | 11 |
| 12672-29-6 | Aroclor-1248 | 8.0 | U | 33 | 8.0 |
| 11097-69-1 | Aroclor-1254 | 14 | U | 33 | 14 |
| 11096-82-5 | Aroclor-1260 | 9.0 | U | 33 | 9.0 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 106 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 110 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 240-212544/8-A

Matrix: Solid | Lab File ID: P1200020.D

Analysis Method: 8082A | Date Collected:

Extraction Method: 3540C | Date Extracted: 12/30/2015 06:54

Sample wt/vol: 30(g) | Date Analyzed: 01/04/2016 13:01

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212845 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 314 | | 33 | 12 |
| 11096-82-5 | Aroclor-1260 | 337 | | 33 | 9.0 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 102 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 110 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-59503-G-4-B MS

Matrix: Solid | Lab File ID: P1200017.D

Analysis Method: 8082A | Date Collected: 12/29/2015 14:14

Extraction Method: 3540C | Date Extracted: 12/30/2015 06:54

Sample wt/vol: 29.79(g) | Date Analyzed: 01/04/2016 12:07

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 24.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212845 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 175 | | 44 | 16 |
| 11104-28-2 | Aroclor-1221 | 21 | U | 44 | 21 |
| 11141-16-5 | Aroclor-1232 | 27 | U | 44 | 27 |
| 53469-21-9 | Aroclor-1242 | 15 | U | 44 | 15 |
| 12672-29-6 | Aroclor-1248 | 11 | U | 44 | 11 |
| 11097-69-1 | Aroclor-1254 | 19 | U | 44 | 19 |
| 11096-82-5 | Aroclor-1260 | 143 | | 44 | 12 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 44 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 43 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-59482-1

SDG No.:

Client Sample ID: | Lab Sample ID: 240-59503-G-4-C MSD

Matrix: Solid | Lab File ID: P1200018.D

Analysis Method: 8082A | Date Collected: 12/29/2015 14:14

Extraction Method: 3540C | Date Extracted: 12/30/2015 06:54

Sample wt/vol: 29.86(g) | Date Analyzed: 01/04/2016 12:25

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-1 ID: 0.53(mm)

% Moisture: 24.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 212845 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 148 | | 44 | 16 |
| 11104-28-2 | Aroclor-1221 | 21 | U | 44 | 21 |
| 11141-16-5 | Aroclor-1232 | 26 | U | 44 | 26 |
| 53469-21-9 | Aroclor-1242 | 15 | U | 44 | 15 |
| 12672-29-6 | Aroclor-1248 | 11 | U | 44 | 11 |
| 11097-69-1 | Aroclor-1254 | 19 | U | 44 | 19 |
| 11096-82-5 | Aroclor-1260 | 118 | | 44 | 12 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 31 | p | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 37 | | 10-155 |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A2HP12 Start Date: 11/16/2015 20:14

Analysis Batch Number: 207007 End Date: 11/17/2015 07:24

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD005 240-207007/4 IC | | 11/16/2015 20:14 | 1 | P1200039.D | CLP-1 0.53(mm) |
| STD005 240-207007/4 IC | | 11/16/2015 20:14 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/5 IC | | 11/16/2015 20:32 | 1 | P1200040.D | CLP-1 0.53(mm) |
| STD01 240-207007/5 IC | | 11/16/2015 20:32 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/6 IC | | 11/16/2015 20:50 | 1 | P1200041.D | CLP-1 0.53(mm) |
| STD02 240-207007/6 IC | | 11/16/2015 20:50 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/7 IC | | 11/16/2015 21:09 | 1 | P1200042.D | CLP-1 0.53(mm) |
| STD05 240-207007/7 IC | | 11/16/2015 21:09 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/8 IC | | 11/16/2015 21:27 | 1 | P1200043.D | CLP-1 0.53(mm) |
| STD1 240-207007/8 IC | | 11/16/2015 21:27 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/9 IC | | 11/16/2015 21:45 | 1 | P1200044.D | CLP-1 0.53(mm) |
| STD2 240-207007/9 IC | | 11/16/2015 21:45 | 1 | | CLP-2 0.53(mm) |
| STD005 240-207007/10 IC | | 11/16/2015 22:03 | 1 | P1200045.D | CLP-1 0.53(mm) |
| STD005 240-207007/10 IC | | 11/16/2015 22:03 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/11 IC | | 11/16/2015 22:21 | 1 | P1200046.D | CLP-1 0.53(mm) |
| STD01 240-207007/11 IC | | 11/16/2015 22:21 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/12 IC | | 11/16/2015 22:39 | 1 | P1200047.D | CLP-1 0.53(mm) |
| STD02 240-207007/12 IC | | 11/16/2015 22:39 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/13 IC | | 11/16/2015 22:57 | 1 | P1200048.D | CLP-1 0.53(mm) |
| STD05 240-207007/13 IC | | 11/16/2015 22:57 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/14 IC | | 11/16/2015 23:15 | 1 | P1200049.D | CLP-1 0.53(mm) |
| STD1 240-207007/14 IC | | 11/16/2015 23:15 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/15 IC | | 11/16/2015 23:33 | 1 | P1200050.D | CLP-1 0.53(mm) |
| STD2 240-207007/15 IC | | 11/16/2015 23:33 | 1 | | CLP-2 0.53(mm) |
| STD005 240-207007/16 IC | | 11/16/2015 23:51 | 1 | P1200051.D | CLP-1 0.53(mm) |
| STD005 240-207007/16 IC | | 11/16/2015 23:51 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/17 IC | | 11/17/2015 00:09 | 1 | P1200052.D | CLP-1 0.53(mm) |
| STD01 240-207007/17 IC | | 11/17/2015 00:09 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/18 IC | | 11/17/2015 00:28 | 1 | P1200053.D | CLP-1 0.53(mm) |
| STD02 240-207007/18 IC | | 11/17/2015 00:28 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/19 IC | | 11/17/2015 00:46 | 1 | P1200054.D | CLP-1 0.53(mm) |
| STD05 240-207007/19 IC | | 11/17/2015 00:46 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/20 IC | | 11/17/2015 01:04 | 1 | P1200055.D | CLP-1 0.53(mm) |

8082A

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A2HP12 Start Date: 11/16/2015 20:14

Analysis Batch Number: 207007 End Date: 11/17/2015 07:24

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-207007/20 IC | | 11/17/2015 01:04 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/21 IC | | 11/17/2015 01:22 | 1 | P1200056.D | CLP-1 0.53(mm) |
| STD2 240-207007/21 IC | | 11/17/2015 01:22 | 1 | | CLP-2 0.53(mm) |
| STD005 240-207007/22 IC | | 11/17/2015 01:40 | 1 | P1200057.D | CLP-1 0.53(mm) |
| STD005 240-207007/22 IC | | 11/17/2015 01:40 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/23 IC | | 11/17/2015 01:58 | 1 | P1200058.D | CLP-1 0.53(mm) |
| STD01 240-207007/23 IC | | 11/17/2015 01:58 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/24 IC | | 11/17/2015 02:16 | 1 | P1200059.D | CLP-1 0.53(mm) |
| STD02 240-207007/24 IC | | 11/17/2015 02:16 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/25 IC | | 11/17/2015 02:34 | 1 | P1200060.D | CLP-1 0.53(mm) |
| STD05 240-207007/25 IC | | 11/17/2015 02:34 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/26 IC | | 11/17/2015 02:52 | 1 | P1200061.D | CLP-1 0.53(mm) |
| STD1 240-207007/26 IC | | 11/17/2015 02:52 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/27 IC | | 11/17/2015 03:10 | 1 | P1200062.D | CLP-1 0.53(mm) |
| STD2 240-207007/27 IC | | 11/17/2015 03:10 | 1 | | CLP-2 0.53(mm) |
| STD005 240-207007/28 IC | | 11/17/2015 03:29 | 1 | P1200063.D | CLP-1 0.53(mm) |
| STD005 240-207007/28 IC | | 11/17/2015 03:29 | 1 | | CLP-2 0.53(mm) |
| STD01 240-207007/29 IC | | 11/17/2015 03:47 | 1 | P1200064.D | CLP-1 0.53(mm) |
| STD01 240-207007/29 IC | | 11/17/2015 03:47 | 1 | | CLP-2 0.53(mm) |
| STD02 240-207007/30 IC | | 11/17/2015 04:05 | 1 | P1200065.D | CLP-1 0.53(mm) |
| STD02 240-207007/30 IC | | 11/17/2015 04:05 | 1 | | CLP-2 0.53(mm) |
| STD05 240-207007/31 ICIS | | 11/17/2015 04:23 | 1 | P1200066.D | CLP-1 0.53(mm) |
| STD05 240-207007/31 ICIS | | 11/17/2015 04:23 | 1 | | CLP-2 0.53(mm) |
| STD1 240-207007/32 IC | | 11/17/2015 04:41 | 1 | P1200067.D | CLP-1 0.53(mm) |
| STD1 240-207007/32 IC | | 11/17/2015 04:41 | 1 | | CLP-2 0.53(mm) |
| STD2 240-207007/33 IC | | 11/17/2015 04:59 | 1 | P1200068.D | CLP-1 0.53(mm) |
| STD2 240-207007/33 IC | | 11/17/2015 04:59 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/34 | | 11/17/2015 05:17 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/34 | | 11/17/2015 05:17 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/35 | | 11/17/2015 05:35 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/35 | | 11/17/2015 05:35 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/36 | | 11/17/2015 05:53 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/36 | | 11/17/2015 05:53 | 1 | | CLP-2 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Instrument ID: A2HP12  Start Date: 11/16/2015 20:14

Analysis Batch Number: 207007  End Date: 11/17/2015 07:24

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| ICV 240-207007/37 | | 11/17/2015 06:12 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/37 | | 11/17/2015 06:12 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/38 | | 11/17/2015 06:30 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/38 | | 11/17/2015 06:30 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/39 | | 11/17/2015 06:48 | 1 | P1200074.D | CLP-1 0.53(mm) |
| ICV 240-207007/39 | | 11/17/2015 06:48 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/40 | | 11/17/2015 07:06 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/40 | | 11/17/2015 07:06 | 1 | | CLP-2 0.53(mm) |
| ICV 240-207007/41 | | 11/17/2015 07:24 | 1 | | CLP-1 0.53(mm) |
| ICV 240-207007/41 | | 11/17/2015 07:24 | 1 | | CLP-2 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A2HP12 Start Date: 01/04/2016 07:42

Analysis Batch Number: 212845 End Date: 01/04/2016 17:15

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-212845/3 CCVIS | | 01/04/2016 07:42 | 1 | P1200003.D | CLP-1 0.53(mm) |
| CCV 240-212845/3 CCVIS | | 01/04/2016 07:42 | 1 | | CLP-2 0.53(mm) |
| CCV 240-212845/4 | | 01/04/2016 08:00 | 1 | P1200004.D | CLP-1 0.53(mm) |
| CCV 240-212845/4 | | 01/04/2016 08:00 | 1 | | CLP-2 0.53(mm) |
| CCV 240-212845/5 | | 01/04/2016 08:19 | 1 | P1200005.D | CLP-1 0.53(mm) |
| CCV 240-212845/5 | | 01/04/2016 08:19 | 1 | | CLP-2 0.53(mm) |
| CCV 240-212845/6 | | 01/04/2016 08:37 | 1 | P1200006.D | CLP-1 0.53(mm) |
| CCV 240-212845/6 | | 01/04/2016 08:37 | 1 | | CLP-2 0.53(mm) |
| CCV 240-212845/7 | | 01/04/2016 08:55 | 1 | P1200007.D | CLP-1 0.53(mm) |
| CCV 240-212845/7 | | 01/04/2016 08:55 | 1 | | CLP-2 0.53(mm) |
| 240-59482-2 | F-VB-UST-P-7 | 01/04/2016 10:19 | 1 | P1200011.D | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 10:19 | 1 | | CLP-2 0.53(mm) |
| 240-59482-3 | F-VB-UST-P-8 | 01/04/2016 10:37 | 1 | P1200012.D | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 10:37 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 10:55 | 20 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 10:55 | 20 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 11:13 | 50 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 11:13 | 50 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 11:49 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 11:49 | 1 | | CLP-2 0.53(mm) |
| 240-59503-G-4-B MS | | 01/04/2016 12:07 | 1 | P1200017.D | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 12:07 | 1 | | CLP-2 0.53(mm) |
| 240-59503-G-4-C MSD | | 01/04/2016 12:25 | 1 | P1200018.D | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 12:25 | 1 | | CLP-2 0.53(mm) |
| MB 240-212544/7-A | | 01/04/2016 12:43 | 1 | P1200019.D | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 12:43 | 1 | | CLP-2 0.53(mm) |
| LCS 240-212544/8-A | | 01/04/2016 13:01 | 1 | P1200020.D | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 13:01 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 13:56 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 13:56 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 14:14 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 14:14 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 14:32 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 14:32 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 14:50 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 14:50 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 15:08 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 15:08 | 1 | | CLP-2 0.53(mm) |
| CCVIS 240-212845/28 | | 01/04/2016 15:26 | 1 | | CLP-1 0.53(mm) |
| CCVIS 240-212845/28 | | 01/04/2016 15:26 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 16:21 | 1 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 16:21 | 1 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 16:39 | 2 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 16:39 | 2 | | CLP-2 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: A2HP12 Start Date: 01/04/2016 07:42

Analysis Batch Number: 212845 End Date: 01/04/2016 17:15

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| ZZZZZ | | 01/04/2016 16:57 | 2 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 16:57 | 2 | | CLP-2 0.53(mm) |
| ZZZZZ | | 01/04/2016 17:15 | 5 | | CLP-1 0.53(mm) |
| ZZZZZ | | 01/04/2016 17:15 | 5 | | CLP-2 0.53(mm) |

# PCBS BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Batch Number: 212544  Batch Start Date: 12/30/15 08:55  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 12/31/15 12:55

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | ex10PPMSPK 00019 | ex2/.2SURRW 00063 | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-59482-2 | F-VB-UST-P-7 | 3540C, 8082A | T | 29.83 g | 10 mL | | 1 mL | | |
| 240-59482-3 | F-VB-UST-P-8 | 3540C, 8082A | T | 29.77 g | 10 mL | | 1 mL | | |
| 240-59503-G-4-B MS | | 3540C, 8082A | T | 29.79 g | 10 mL | 1 mL | 1 mL | | |
| 240-59503-G-4-C MSD | | 3540C, 8082A | T | 29.86 g | 10 mL | 1 mL | 1 mL | | |
| MB 240-212544/7 | | 3540C, 8082A | | 30 g | 10 mL | | 1 mL | | |
| LCS 240-212544/8 | | 3540C, 8082A | | 30 g | 10 mL | 1 mL | 1 mL | | |

| Batch Notes | |
|---|---|
| Acid used for Clean Up Reagent | 2226509 |
| Balance ID | 1339911 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2345798 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GENERAL CHEMISTRY

COVER PAGE
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job Number: 240-59482-1

SDG No.:

Project: MRC Block F UST Pipe Inv.

| Client Sample ID | Lab Sample ID |
|---|---|
| F-VB-UST-P-7 | 240-59482-2 |
| F-VB-UST-P-8 | 240-59482-3 |

Comments:

9-IN
DETECTION LIMITS
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job Number: 240-59482-1

SDG Number:

Matrix: Solid Instrument ID: NOEQUIP

Method: Moisture RL Date: 01/28/2010 09:24

| Analyte | Wavelength/ Mass | RL (%) | |
|---|---|---|---|
| Percent Moisture | | 0.1 | |
| Percent Solids | | 0.1 | |

9-IN
CALIBRATION BLANK DETECTION LIMITS
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton  Job Number: 240-59482-1

SDG Number:

Matrix: Solid  Instrument ID: NOEQUIP

Method: Moisture  XRL Date: 01/28/2010 09:24

| Analyte | Wavelength/ Mass | XRL (mg/L) | |
|---|---|---|---|
| Percent Moisture | | 10 | |
| Percent Solids | | 10 | |

13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Canton Job No.: 240-59482-1

SDG No.:

Instrument ID: NOEQUIP Method: Moisture

Start Date: 12/29/2015 13:50 End Date: 12/29/2015 13:50

| Lab Sample ID | D / F | Type | Time | Analytes | | | | | | | | | | | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | | | | % Sol | Moist | | | | | | | | | | | | | | | | | | |
| 240-59255-B-13 DU | 1 | T | 13:50 | X | X | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| 240-59482-2 | 1 | T | 13:50 | X | X | | | | | | | | | | | | | | | | | | |
| 240-59482-3 | 1 | T | 13:50 | X | X | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |
| ZZZZZZ | | | 13:50 | | | | | | | | | | | | | | | | | | | | |

Prep Types
T = Total/NA

# GENERAL CHEMISTRY BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-59482-1

SDG No.:

Batch Number: 212487  Batch Start Date: 12/29/15 13:50  Batch Analyst: Woodward, Bruce

Batch Method: Moisture  Batch End Date: 12/30/15 08:40

| Lab Sample ID | Client Sample ID | Method Chain | Basis | DishWeight | SampleMassWet | SampleMassDry | | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-59255-B-13 DU | | Moisture | T | 4.3859 g | 9.1871 g | 9.1038 g | | | |
| 240-59482-2 | F-VB-UST-P-7 | Moisture | T | 4.3859 g | 21.6950 g | 18.5049 g | | | |
| 240-59482-3 | F-VB-UST-P-8 | Moisture | T | 4.3859 g | 19.3165 g | 16.9117 g | | | |

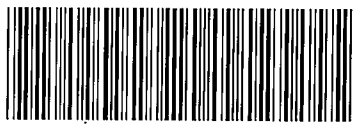
| Batch Notes | |
|---|---|
| Balance ID | B47 No Unit |
| Date and Time Samples in Desiccator | 12/30/15; 04:36 |
| Date and Time Samples out of Desiccator | 12/30/15; 8:38 |
| Date samples were placed in the oven | 12/29/15 |
| Oven Temp when samples are put in oven | 103.1 Degrees C |
| Time samples were place in the oven | 13:55 |
| Date samples were removed from oven | 12/30/15 |
| Oven Temp when samples removed from oven | 103.3 Degrees C |
| Time Samples were removed from oven | 04:36 |
| Oven ID | 002 |
| ID number of the thermometer | TempGuard Box-C #6 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# Shipping and Receiving Documents

# CHAIN OF CUSTODY
# AND
# RECEIVING DOCUMENTS



240-59482 Chain of Custody



4101 Shuffel Street, N.W. North Canton, OH 44720 **tel** 330.497.9396 **fax** 330.497.0772 **www.testamericainc.com**

2.8/c 2.5

# Chain of Custody Record

016118

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING
TestAmerica Laboratories, Inc.
TAL-8210 (0713)

Regulatory Program: ☐ DW ☐ NPDES ☐ RCRA ☐ Other:

| Client Contact | Project Manager: SCOTT NESBIT | Site Contact: MULLIS | Date: 12/28/15 | COC No: |
|---|---|---|---|---|
| Company Name: TETRA TECH | Tel/Fax: 412 921 7134 | Lab Contact: MCFADDEN | Carrier: Fed Ex | 1 of 1 COCs |
| Address: 20251 CENTURY BLVD. STE.200 | Analysis Turnaround Time | | | Sampler: |
| City/State/Zip: GERMANTOWN, MD, 20874 | ☐ CALENDAR DAYS ☐ WORKING DAYS | | | For Lab Use Only: |
| Phone: 301-528-3000 | TAT if different from Below | | | Walk-in Client: |
| Fax: 301-528-3000 | ☐ 2 weeks ☐ 1 week (5 DAYS) ☐ 2 days ☐ 1 day | | | Lab Sampling: |
| Project Name: MRC SOIL REMEDY, Block F | | | | Job / SDG No.: |
| Site: Block F UST PIPE INV. | | | | |
| P O # 112IC07336 | | | | |

| Sample Identification | Sample Date | Sample Time | Sample Type (C=Comp, G=Grab) | Matrix | # of Cont. | Filtered Sample (Y/N) | Perform MS / MSD (Y/N) | VOCs (8260) | TPH-GRO/DRO (8015) | PCBs (8082) | SVOCs (8270) | Sample Specific Notes: |
|---|---|---|---|---|---|---|---|---|---|---|---|---|
| TB-122815 | 12/28/15 | — | G | AQ | 2 | | | X | | | | |
| F-VB-UST-P-7 | 1 | 1000 | 1 | SO | 3 | | | X | X | X | X | |
| F-VB-UST-P-8 | 1 | 1030 | 1 | SO | 3 | | | X | X | X | X | |

12/28/15

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Preservation Used: 1= Ice, 2= HCl; 3= H2SO4; 4=HNO3; 5=NaOH; 6= Other _____

**Possible Hazard Identification:**
Are any samples from a listed EPA Hazardous Waste? Please List any EPA Waste Codes for the sample in the Comments Section if the lab is to dispose of the sample.

☑ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

**Sample Disposal ( A fee may be assessed if samples are retained longer than 1 month)**

☐ Return to Client ☑ Disposal by Lab ☐ Archive for _____ Months

**Special Instructions/QC Requirements & Comments:** MRC SOIL REMEDY; Block F UST PIPE REMOVAL INV., BASE SAMPLES

| Custody Seals Intact: ☐ Yes ☐ No | Custody Seal No.: | | Cooler Temp. (°C): Obs'd: _____ Corr'd: _____ Therm ID No.: _____ | | |
|---|---|---|---|---|---|
| Relinquished by: | Company: Tetra Tech | Date/Time: 12/28/15 | Received by: | Company: | Date/Time: |
| Relinquished by: | Company: | Date/Time: | Received by: | Company: | Date/Time: |
| Relinquished by: | Company: | Date/Time: | Received in Laboratory by: Jacquelyn Reed | Company: TANC | Date/Time: 12/29/15 930 |

TestAmerica Canton Sample Receipt Form/Narrative
Canton Facility

Login # : 594482

Client Tetra Tech Site Name

Cooler unpacked by: A. Jerel

Cooler Received on 12/29/15 Opened on 12/29/15

FedEx: 1st Grd Exp UPS FAS Stetson Client Drop Off TestAmerica Courier Other

Receipt After-hours: Drop-off Date/Time Storage Location

TestAmerica Cooler # Foam Box Client Cooler Box Other

Packing material used: Bubble Wrap Foam Plastic Bag None Other

COOLANT: Wet Ice Blue Ice Dry Ice Water None

1. Cooler temperature upon receipt

IR GUN# 53 (CF +0.1 °C) Observed Cooler Temp. °C Corrected Cooler Temp. °C

IR GUN# 48 (CF -0.3 °C) Observed Cooler Temp. 2.8 °C Corrected Cooler Temp. 2.5 °C ☐ See Multiple Cooler Form

~~IR GUN# 5 (CF +0.4 °C) Observed Cooler Temp. °C Corrected Cooler Temp. °C~~

IR GUN# 8 (CF -0.5 °C) Observed Cooler Temp. °C Corrected Cooler Temp. °C

| | | |
|---|---|---|
| 2. Were custody seals on the outside of the cooler(s)? If Yes Quantity 1 | Yes | No |
| -Were custody seals on the outside of the cooler(s) signed & dated? | Yes | No NA |
| -Were custody seals on the bottle(s) or bottle kits (LLHg/MeHg)? | Yes | No |
| 3. Shippers' packing slip attached to the cooler(s)? | Yes | No |
| 4. Did custody papers accompany the sample(s)? | Yes | No |
| 5. Were the custody papers relinquished & signed in the appropriate place? | Yes | No |
| 6. Was/were the person(s) who collected the samples clearly identified on the COC? | Yes | No |
| 7. Did all bottles arrive in good condition (Unbroken)? | Yes | No |
| 8. Could all bottle labels be reconciled with the COC? | Yes | No |
| 9. Were correct bottle(s) used for the test(s) indicated? | Yes | No |
| 10. Sufficient quantity received to perform indicated analyses? | Yes | No |
| 11. Were sample(s) at the correct pH upon receipt? | Yes | No NA pH Strip Lot# HC559158 |
| 12. Were VOAs on the COC? | Yes | No |
| 13. Were air bubbles >6 mm in any VOA vials? | Yes | No NA |
| 14. Was a VOA trip blank present in the cooler(s)? Trip Blank Lot # N/A | Yes | No |
| 15. Was a LL Hg or Me Hg trip blank present? | Yes | No |

Contacted PM Date by via Verbal Voice Mail Other

Concerning

14. CHAIN OF CUSTODY & SAMPLE DISCREPANCIES

Samples processed by:

15. SAMPLE CONDITION

Sample(s) were received after the recommended holding time had expired.

Sample(s) were received in a broken container.

Sample(s) were received with bubble >6 mm in diameter. (Notify PM)

16. SAMPLE PRESERVATION

Sample(s) were further preserved in the laboratory.

Time preserved: Preservative(s) added/Lot number(s):

# VALIDATION REPORTS

TETRA TECH

# INTERNAL CORRESPONDENCE

**TO:** **S. NESBIT** **DATE:** **NOVEMBER 25, 2015**

**FROM:** **MICHELLE L. ALLEN** **COPIES:** **DV FILE**

**SUBJECT:** **ORGANIC & INORGANIC DATA VALIDATION – VOC/PAH/PCB/DRO/GRO**
**LOCKHEED MARTIN CORPORATION (LMC) – BLOCK F**
**UNDERGROUND STORAGE TANK (UST) AREA**
**SDG 240-57597-1**

**SAMPLES:** 15/Soil/VOC/PAH/PCB/DRO/GRO/Metals

| | | |
|---|---|---|
| F-VB-UST-1-14 | F-VB-UST-10-14 | F-VB-UST-11-14 |
| F-VB-UST-12-14 | F-VB-UST-13-14 | F-VB-UST-14-14 |
| F-VB-UST-2-14 | F-VB-UST-3-14 | F-VB-UST-4-14 |
| F-VB-UST-5-14 | F-VB-UST-6-14 | F-VB-UST-7-14 |
| F-VB-UST-8-14 | F-VB-UST-9-14 | F-VB-UST-DUP |

1/Aqueous/VOC

TB-110515

Overview

The sample set for LMC Block F UST Area, SDG 240-57597-1 consisted of fifteen (15) soil environmental samples and one (1) trip blank. All fifteen (15) soil samples were analyzed for Volatile Organic Compounds (VOC), Semi-Volatile Organic Compounds (SVOC), Polychlorinated Biphenyls (PCB), Diesel Range Organics (DRO), and Gasoline Range Organics (GRO). The trip blank was analyzed for VOC only. One field duplicate sample pair was included in this SDG F-VB-UST-2-14/F-VB-UST-DUP.

The samples were collected by Tetra Tech, Inc. on November 5, 2015 and analyzed by TestAmerica, Inc. All analyses were conducted in accordance with SW-846 Methods 8260C, 8270D, 8082A, and 8015C analytical and reporting protocols.

The data contained in this SDG were validated with regard to the following parameters: data completeness, holding times and sample preservation, GC/MS tuning, initial/continuing calibrations, laboratory method/preparation/trip blanks, surrogate spike recoveries, laboratory control sample results, matrix spike/matrix spike duplicate results, internal standard areas, field duplicate precision, chromatographic resolution, compound identification, compound quantitation, and detection limits. Areas of concern are listed below.

Major

No major issues were identified.

Minor

- The continuing calibration performed on instrument A3UX14 on 11/10/15 @ 10:47 had %Ds for trichlorofluoromethane, 1,1,2-trichlorotrifluoromethane, 2,2-dichloropropane, 1,1,1-trichloroethane, and carbon tetrachloride which exceeded the 20% quality control limit. All soil samples, with the exception of sample F-VB-UST-DUP, were affected. The non-detected results reported for these compounds in the affected samples were qualified as estimated, (UJ).

- The continuing calibration performed on instrument A3UX14 on 11/11/15 @ 10:39 had a Percent Drift (%Drift) for acetone and %Ds for tert-butyl alcohol, 2,2-dichloropropane, 2-butanone, 2-

chloroethyl-ethyl vinyl ether, 4-methyl-2-pentanone, and 2-hexanone which exceeded the 20% quality control limit. Sample F-VB-UST-DUP was affected. No action was taken for 2-butanone because this result was qualified for laboratory blank contamination. The non-detected results reported for the remaining compounds in the affected sample were qualified as estimated, (UJ).

- The continuing calibration performed on instrument A3UX15 on 11/12/15 @ 10:41 had a %Ds for hexachlorobutadiene which exceeded the 20% quality control limit. The non-detected result reported for this compound in the affected sample, TB-110515, was qualified as estimated, (UJ).

- The continuing calibration performed on instrument A4HP10 on 11/11/15 @ 09:34 had %Ds for benzaldehyde and hexachlorocyclopentadiene which exceeded the 20% quality control limit. Samples F-VB-UST-1-14, F-VB-UST-2-14, F-VB-UST-3-14, F-VB-UST-4-14, F-VB-UST-5-14, F-VB-UST-6-14, F-VB-UST-7-14, F-VB-UST-8-14, F-VB-UST-9-14, and F-VB-UST-10-14 were affected. The non-detected results reported for these compounds in the affected samples were qualified as estimated, (UJ).

- The continuing calibration performed on instrument A4HP9 on 11/11/15 @ 07:49 had %Ds for benzaldehyde and N-nitroso-di-n-propylamine which exceeded the 20% quality control limit. Samples F-VB-UST-11-14, F-VB-UST-12-14, F-VB-UST-13-14, F-VB-UST-14-14, and F-VB-UST-DUP were affected. The detected and non-detected results reported for these compounds in the affected samples were qualified as estimated, (J) and (UJ), respectively.

- The following compounds were detected in the laboratory method and/or trip blanks at the following maximum concentrations:

| Compound | Maximum Concentration | Action Level |
|---|---|---|
| 2-Butanone (1) | 3.14 µg/Kg | 31.4 µg/Kg |
| Naphthalene (1) | 0.35 µg/Kg | 1.75 µg/Kg |
| Toluene (1) | 0.369 µg/Kg | 1.85 µg/Kg |
| 2-Butanone (2) | 2.83 µg/Kg | 28.3 µg/Kg |
| Tert-butyl alcohol (2) | 6.1 µg/Kg | 30.5 µg/Kg |
| DRO (3) | 7.82 mg/Kg | 39.1 mg/Kg |

(1) - Maximum concentration detected in the laboratory method blank, MB 240-206267/6, affecting sample F-VB-UST-DUP from preparation batch # 240-206267.
(2) - Maximum concentration detected in the laboratory method blank, MB 240-206089/1-A, affecting all samples, except sample F-VB-UST-DUP, from preparation batch # 240-206089.
(3) - Maximum concentration detected in the laboratory method blank, MB 240-205788/22-A, affecting all samples.

An action level of 10X the maximum concentration for the common laboratory contaminant, 2-butanone, and 5X for the remaining compounds were established to evaluate for blank contamination. Percent moisture and dilution factor, if necessary, were taken into consideration during the application of the action level. The detected results reported for these compounds in the affected samples below the established action level were qualified as non-detected, (U).

- Detected results reported below the RL but above the Method Detection Limit (MDL) were qualified as estimated, (J).

Notes

Toxicity Characteristic Leaching Procedure (TCLP) metals were requested on the sample Chains of Custody (COCs), however, the client cancelled this analysis request for all the samples.

2-Butanone was detected in the laboratory method blank associated with the trip blank. No action was taken because field blanks are not qualified for laboratory blank contamination.

The VOC Matrix Spike/Matrix Spike Duplicate (MS/MSD) analyses performed on sample F-VB-UST-5-14 had Percent Recoveries (%Rs) for 4-methyl-2-pentanone above the upper quality control limit. In addition, the MS %R was low for vinyl acetate. The Relative Percent Differences (RPDs) for hexachlorobutadiene, n-butylbenzene, sec-butylbenzene, and 4-isopropyltoluene exceeded the 30% quality control criterion. No validation action was taken because 4-methyl-2-pentanone was not detected in the parent sample, the MSD %R for vinyl acetate was acceptable, and the MS/MSD %Rs for the remaining compounds were within the quality control limits.

The GRO MS/MSD analyses performed on F-VB-UST-DUP had a low MSD %R and the RPD exceeded the quality control limit. No action was taken because the MS and associated Laboratory Control Sample (LCS) %Rs were acceptable.

Naphthalene was reported in both the VOC and PAH fractions.

In the PCB fraction, samples F-VB-UST-5-14 and F-VB-UST-13-14 required a tetrabutylammonium sulfite (TBA) clean-up to reduce matrix interferences caused by sulfur.

Non-detected results were reported to the MDL.

**Executive Summary**

**Laboratory Performance:** VOC and SVOC continuing calibration %Ds or %Drift exceeded 20%. Contaminants were detected in the VOC and DROs laboratory method blanks.

**Other Factors Affecting Data Quality:** The VOC MS/MSD analyses had noncompliant %Rs and RPDs. The GRO MSD %R was low. Results below the RL were estimated.

The data for these analyses were reviewed with reference to the "National Functional Guidelines for Organic Review" (June 2008) and the "National Functional Guidelines for Inorganic Review" (January 2010). The text of this report has been formulated to address only those areas affecting data quality.

Tetra Tech, Inc.
Michelle L. Allen
Chemist/Data Validator

Tetra Tech, Inc.
Joseph A. Samchuck
Data Validation Manager

Attachments:
Appendix A – Qualified Analytical Results
Appendix B – Results as Reported by the Laboratory
Appendix C – Support Documentation

**Appendix A**

Qualified Analytical Results

**Qualifier Codes:**

A = Lab Blank Contamination
B = Field Blank Contamination
C = Calibration Noncompliance (i.e., % RSDs, %Ds, ICVs, CCVs, RRFs, etc.)
C01 = GC/MS Tuning Noncompliance
D = MS/MSD Recovery Noncompliance
E = LCS/LCSD Recovery Noncompliance
F = Lab Duplicate Imprecision
G = Field Duplicate Imprecision
H = Holding Time Exceedance
I = ICP Serial Dilution Noncompliance
J = ICP PDS Recovery Noncompliance; MSA's $r < 0.995$
K = ICP Interference - includes ICS % R Noncompliance
L = Instrument Calibration Range Exceedance
M = Sample Preservation Noncompliance
N = Internal Standard Noncompliance
N01 = Internal Standard Recovery Noncompliance Dioxins
N02 = Recovery Standard Noncompliance Dioxins
N03 = Clean-up Standard Noncompliance Dioxins
O = Poor Instrument Performance (i.e., base-time drifting)
P = Uncertainty near detection limit (< 2 x IDL for inorganics and <CRQL for organics)
Q = Other problems (can encompass a number of issues; i.e.chromatography,interferences, etc.)
R = Surrogates Recovery Noncompliance
S = Pesticide/PCB Resolution
T = % Breakdown Noncompliance for DDT and Endrin
U = RPD between columns/detectors >40% for positive results determined via GC/HPLC
V = Non-linear calibrations; correlation coefficient $r < 0.995$
W = EMPC result
X = Signal to noise response drop
Y = Percent solids <30%
Z = Uncertainty at 2 standard deviations is greater than sample activity
Z1 = Tentatively Identified Compound considered presumptively present
Z2 = Tentatively Identified Compound column bleed
Z3 = Tentatively Identified Compound aldol condensate
Z4 = Sample activity is less than the at uncertainty at 3 standard deviations and greater than the MDC
Z5 = Sample activity is less than the at uncertainty at 3 standard deviations and less than the MDC

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-10-14 | | | F-VB-UST-11-14 | | | F-VB-UST-1-14 | | | F-VB-UST-12-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-10 | | | 240-57597-11 | | | 240-57597-1 | | | 240-57597-12 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 86.6 | | | 83.5 | | | 84.3 | | | 84.1 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| 1,1,1,2-TETRACHLOROETHANE | | 0.56 | U | | 0.6 | U | | 0.5 | U | | 0.55 | U | |
| 1,1,1-TRICHLOROETHANE | | 0.73 | UJ | C | 0.77 | UJ | C | 0.65 | UJ | C | 0.72 | UJ | C |
| 1,1,2,2-TETRACHLOROETHANE | | 0.35 | U | | 0.37 | U | | 0.31 | U | | 0.34 | U | |
| 1,1,2-TRICHLOROTRIFLUOROETHANE | | 0.47 | UJ | C | 0.5 | UJ | C | 0.42 | UJ | C | 0.46 | UJ | C |
| 1,1-DICHLOROETHANE | | 0.38 | U | | 0.4 | U | | 0.34 | U | | 0.38 | U | |
| 1,1-DICHLOROETHENE | | 0.9 | U | | 0.95 | U | | 0.81 | U | | 0.89 | U | |
| 1,1-DICHLOROPROPENE | | 0.42 | U | | 0.44 | U | | 0.37 | U | | 0.41 | U | |
| 1,2,3-TRICHLOROBENZENE | | 0.56 | U | | 0.6 | U | | 0.5 | U | | 0.55 | U | |
| 1,2,3-TRICHLOROPROPANE | | 0.88 | U | | 0.93 | U | | 0.79 | U | | 0.86 | U | |
| 1,2,3-TRIMETHYLBENZENE | | 0.52 | U | | 0.55 | U | | 0.46 | U | | 0.51 | U | |
| 1,2,4-TRICHLOROBENZENE | | 0.43 | U | | 0.45 | U | | 0.38 | U | | 0.42 | U | |
| 1,2,4-TRIMETHYLBENZENE | | 0.65 | U | | 0.69 | U | | 0.58 | U | | 0.64 | U | |
| 1,2-DIBROMO-3-CHLOROPROPANE | | 3 | U | | 3.2 | U | | 2.7 | U | | 3 | U | |
| 1,2-DIBROMOETHANE | | 0.6 | U | | 0.63 | U | | 0.53 | U | | 0.59 | U | |
| 1,2-DICHLOROBENZENE | | 0.38 | U | | 0.4 | U | | 0.34 | U | | 0.38 | U | |
| 1,2-DICHLOROETHANE | | 0.52 | U | | 0.55 | U | | 0.46 | U | | 0.51 | U | |
| 1,2-DICHLOROPROPANE | | 0.16 | U | | 0.17 | U | | 0.14 | U | | 0.15 | U | |
| 1,3-DICHLOROBENZENE | | 0.78 | U | | 0.82 | U | | 0.69 | U | | 0.76 | U | |
| 1,3-DICHLOROPROPANE | | 0.45 | U | | 0.48 | U | | 0.4 | U | | 0.44 | U | |
| 1,4-DICHLOROBENZENE | | 0.31 | U | | 0.33 | U | | 0.28 | U | | 0.31 | U | |
| 2,2-DICHLOROPROPANE | | 0.49 | UJ | C | 0.52 | UJ | C | 0.44 | UJ | C | 0.49 | UJ | C |
| 2-BUTANONE | | 3.5 | U | A | 3.9 | U | A | 3.9 | U | A | 3.7 | U | A |
| 2-CHLOROETHYL VINYL ETHER | | 0.72 | U | | 0.76 | U | | 0.64 | U | | 0.71 | U | |
| 2-CHLOROTOLUENE | | 0.34 | U | | 0.36 | U | | 0.3 | U | | 0.33 | U | |
| 2-HEXANONE | | 0.76 | U | | 0.81 | U | | 0.68 | U | | 0.75 | U | |
| 4-CHLOROTOLUENE | | 0.29 | U | | 0.31 | U | | 0.26 | U | | 0.29 | U | |
| 4-ISOPROPYLTOLUENE | | 0.46 | U | | 0.49 | U | | 0.41 | U | | 0.45 | U | |
| 4-METHYL-2-PENTANONE | | 1.4 | U | | 1.5 | U | | 1.2 | U | | 1.4 | U | |
| ACETONE | | 6.4 | U | | 6.8 | U | | 5.7 | U | | 6.3 | U | |
| BENZENE | | 1 | U | | 1.1 | U | | 0.91 | U | | 1 | U | |
| BROMOBENZENE | | 0.37 | U | | 0.39 | U | | 0.33 | U | | 0.37 | U | |
| BROMOCHLOROMETHANE | | 0.37 | U | | 0.39 | U | | 0.33 | U | | 0.37 | U | |
| BROMODICHLOROMETHANE | | 0.33 | U | | 0.35 | U | | 0.29 | U | | 0.32 | U | |
| BROMOFORM | | 0.27 | U | | 0.29 | U | | 0.24 | U | | 0.27 | U | |
| BROMOMETHANE | | 0.47 | U | | 0.5 | U | | 0.42 | U | | 0.46 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-13-14 | | | F-VB-UST-14-14 | | | F-VB-UST-2-14 | | | F-VB-UST-3-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-13 | | | 240-57597-14 | | | 240-57597-2 | | | 240-57597-3 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 85.3 | | | 85.5 | | | 85.0 | | | 82.2 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| 1,1,1,2-TETRACHLOROETHANE | | 0.52 | U | | 0.53 | U | | 0.55 | U | | 0.58 | U | |
| 1,1,1-TRICHLOROETHANE | | 0.67 | UJ | C | 0.68 | UJ | C | 0.71 | UJ | C | 0.76 | UJ | C |
| 1,1,2,2-TETRACHLOROETHANE | | 0.32 | U | | 0.33 | U | | 0.34 | U | | 0.36 | U | |
| 1,1,2-TRICHLOROTRIFLUOROETHANE | | 0.43 | UJ | C | 0.44 | UJ | C | 0.46 | UJ | C | 0.49 | UJ | C |
| 1,1-DICHLOROETHANE | | 0.35 | U | | 0.36 | U | | 0.37 | U | | 0.4 | U | |
| 1,1-DICHLOROETHENE | | 0.83 | U | | 0.84 | U | | 0.88 | U | | 0.93 | U | |
| 1,1-DICHLOROPROPENE | | 0.38 | U | | 0.39 | U | | 0.41 | U | | 0.43 | U | |
| 1,2,3-TRICHLOROBENZENE | | 0.52 | U | | 0.53 | U | | 0.55 | U | | 0.58 | U | |
| 1,2,3-TRICHLOROPROPANE | | 0.81 | U | | 0.82 | U | | 0.86 | U | | 0.91 | U | |
| 1,2,3-TRIMETHYLBENZENE | | 0.48 | U | | 0.48 | U | | 0.5 | U | | 0.53 | U | |
| 1,2,4-TRICHLOROBENZENE | | 0.39 | U | | 0.4 | U | | 0.42 | U | | 0.44 | U | |
| 1,2,4-TRIMETHYLBENZENE | | 0.6 | U | | 0.61 | U | | 0.64 | U | | 0.67 | U | |
| 1,2-DIBROMO-3-CHLOROPROPANE | | 2.8 | U | | 2.9 | U | | 3 | U | | 3.2 | U | |
| 1,2-DIBROMOETHANE | | 0.55 | U | | 0.56 | U | | 0.58 | U | | 0.62 | U | |
| 1,2-DICHLOROBENZENE | | 0.35 | U | | 0.36 | U | | 0.37 | U | | 0.4 | U | |
| 1,2-DICHLOROETHANE | | 0.48 | U | | 0.48 | U | | 0.5 | U | | 0.53 | U | |
| 1,2-DICHLOROPROPANE | | 0.14 | U | | 0.15 | U | | 0.15 | U | | 0.16 | U | |
| 1,3-DICHLOROBENZENE | | 0.71 | U | | 0.73 | U | | 0.76 | U | | 0.8 | U | |
| 1,3-DICHLOROPROPANE | | 0.41 | U | | 0.42 | U | | 0.44 | U | | 0.46 | U | |
| 1,4-DICHLOROBENZENE | | 0.29 | U | | 0.3 | U | | 0.31 | U | | 0.33 | U | |
| 2,2-DICHLOROPROPANE | | 0.45 | UJ | C | 0.46 | UJ | C | 0.48 | UJ | C | 0.51 | UJ | C |
| 2-BUTANONE | | 3.1 | U | A | 4 | U | A | 4 | U | A | 3.9 | U | A |
| 2-CHLOROETHYL VINYL ETHER | | 0.66 | U | | 0.67 | U | | 0.7 | U | | 0.74 | U | |
| 2-CHLOROTOLUENE | | 0.31 | U | | 0.32 | U | | 0.33 | U | | 0.35 | U | |
| 2-HEXANONE | | 0.7 | U | | 0.72 | U | | 0.75 | U | | 0.79 | U | |
| 4-CHLOROTOLUENE | | 0.27 | U | | 0.27 | U | | 0.29 | U | | 0.3 | U | |
| 4-ISOPROPYLTOLUENE | | 0.42 | U | | 0.43 | U | | 0.45 | U | | 0.48 | U | |
| 4-METHYL-2-PENTANONE | | 1.3 | U | | 1.3 | U | | 1.3 | U | | 1.4 | U | |
| ACETONE | | 5.9 | U | | 6 | U | | 6.3 | U | | 6.6 | U | |
| BENZENE | | 0.93 | U | | 0.95 | U | | 0.99 | U | | 1 | U | |
| BROMOBENZENE | | 0.34 | U | | 0.35 | U | | 0.36 | U | | 0.38 | U | |
| BROMOCHLOROMETHANE | | 0.34 | U | | 0.35 | U | | 0.36 | U | | 0.38 | U | |
| BROMODICHLOROMETHANE | | 0.3 | U | | 0.31 | U | | 0.32 | U | | 0.34 | U | |
| BROMOFORM | | 0.25 | U | | 0.25 | U | | 0.26 | U | | 0.28 | U | |
| BROMOMETHANE | | 0.43 | U | | 0.44 | U | | 0.46 | U | | 0.49 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-4-14 | | | F-VB-UST-5-14 | | | F-VB-UST-6-14 | | | F-VB-UST-7-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-4 | | | 240-57597-5 | | | 240-57597-6 | | | 240-57597-7 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 84.4 | | | 84.8 | | | 83.5 | | | 83.8 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| 1,1,1,2-TETRACHLOROETHANE | | 0.55 | U | | 0.58 | U | | 0.58 | U | | 0.56 | U | |
| 1,1,1-TRICHLOROETHANE | | 0.71 | UJ | C | 0.75 | UJ | C | 0.75 | UJ | C | 0.73 | UJ | C |
| 1,1,2,2-TETRACHLOROETHANE | | 0.34 | U | | 0.36 | U | | 0.36 | U | | 0.35 | U | |
| 1,1,2-TRICHLOROTRIFLUOROETHANE | | 0.46 | UJ | C | 0.49 | UJ | C | 0.49 | UJ | C | 0.47 | UJ | C |
| 1,1-DICHLOROETHANE | | 0.37 | U | | 0.39 | U | | 0.39 | U | | 0.38 | U | |
| 1,1-DICHLOROETHENE | | 0.87 | U | | 0.93 | U | | 0.93 | U | | 0.9 | U | |
| 1,1-DICHLOROPROPENE | | 0.4 | U | | 0.43 | U | | 0.43 | U | | 0.42 | U | |
| 1,2,3-TRICHLOROBENZENE | | 0.55 | U | | 0.58 | U | | 0.58 | U | | 0.56 | U | |
| 1,2,3-TRICHLOROPROPANE | | 0.85 | U | | 0.9 | U | | 0.9 | U | | 0.88 | U | |
| 1,2,3-TRIMETHYLBENZENE | | 0.5 | U | | 0.53 | U | | 0.53 | U | | 0.52 | U | |
| 1,2,4-TRICHLOROBENZENE | | 0.41 | U | | 0.44 | U | | 0.44 | U | | 0.43 | U | |
| 1,2,4-TRIMETHYLBENZENE | | 0.63 | U | | 0.67 | U | | 0.67 | U | | 0.65 | U | |
| 1,2-DIBROMO-3-CHLOROPROPANE | | 3 | U | | 3.1 | U | | 3.1 | U | | 3 | U | |
| 1,2-DIBROMOETHANE | | 0.58 | U | | 0.61 | U | | 0.61 | U | | 0.6 | U | |
| 1,2-DICHLOROBENZENE | | 0.37 | U | | 0.39 | U | | 0.39 | U | | 0.38 | U | |
| 1,2-DICHLOROETHANE | | 0.5 | U | | 0.53 | U | | 0.53 | U | | 0.52 | U | |
| 1,2-DICHLOROPROPANE | | 0.15 | U | | 0.16 | U | | 0.16 | U | | 0.16 | U | |
| 1,3-DICHLOROBENZENE | | 0.75 | U | | 0.8 | U | | 0.8 | U | | 0.78 | U | |
| 1,3-DICHLOROPROPANE | | 0.44 | U | | 0.46 | U | | 0.46 | U | | 0.45 | U | |
| 1,4-DICHLOROBENZENE | | 0.31 | U | | 0.32 | U | | 0.32 | U | | 0.31 | U | |
| 2,2-DICHLOROPROPANE | | 0.48 | UJ | C | 0.51 | UJ | C | 0.51 | UJ | C | 0.49 | UJ | C |
| 2-BUTANONE | | 4 | U | A | 4.3 | U | A | 3.8 | U | A | 4.1 | U | A |
| 2-CHLOROETHYL VINYL ETHER | | 0.7 | U | | 0.74 | U | | 0.74 | U | | 0.72 | U | |
| 2-CHLOROTOLUENE | | 0.33 | U | | 0.35 | U | | 0.35 | U | | 0.34 | U | |
| 2-HEXANONE | | 0.74 | U | | 0.79 | U | | 0.79 | U | | 0.76 | U | |
| 4-CHLOROTOLUENE | | 0.28 | U | | 0.3 | U | | 0.3 | U | | 0.29 | U | |
| 4-ISOPROPYLTOLUENE | | 0.45 | U | | 0.47 | U | | 0.47 | U | | 0.46 | U | |
| 4-METHYL-2-PENTANONE | | 1.3 | U | | 1.4 | U | | 1.4 | U | | 1.4 | U | |
| ACETONE | | 6.2 | U | | 13 | J | P | 6.6 | U | | 6.4 | U | |
| BENZENE | | 0.98 | U | | 1 | U | | 1 | U | | 1 | U | |
| BROMOBENZENE | | 0.36 | U | | 0.38 | U | | 0.38 | U | | 0.37 | U | |
| BROMOCHLOROMETHANE | | 0.36 | U | | 0.38 | U | | 0.38 | U | | 0.37 | U | |
| BROMODICHLOROMETHANE | | 0.32 | U | | 0.34 | U | | 0.34 | U | | 0.33 | U | |
| BROMOFORM | | 0.26 | U | | 0.28 | U | | 0.28 | U | | 0.27 | U | |
| BROMOMETHANE | | 0.46 | U | | 0.49 | U | | 0.49 | U | | 0.47 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-8-14 | | | F-VB-UST-9-14 | | | F-VB-UST-DUP | | |
|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-8 | | | 240-57597-9 | | | 240-57597-15 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 82.0 | | | 80.9 | | | 85.5 | | |
| | DUP_OF | | | | | | | F-VB-UST-2-14 | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| 1,1,1,2-TETRACHLOROETHANE | | 0.58 | U | | 0.58 | U | | 0.53 | U | |
| 1,1,1-TRICHLOROETHANE | | 0.75 | UJ | C | 0.75 | UJ | C | 0.68 | U | |
| 1,1,2,2-TETRACHLOROETHANE | | 0.36 | U | | 0.36 | U | | 0.33 | U | |
| 1,1,2-TRICHLOROTRIFLUOROETHANE | | 0.48 | UJ | C | 0.49 | UJ | C | 0.44 | U | |
| 1,1-DICHLOROETHANE | | 0.39 | U | | 0.39 | U | | 0.36 | U | |
| 1,1-DICHLOROETHENE | | 0.92 | U | | 0.92 | U | | 0.84 | U | |
| 1,1-DICHLOROPROPENE | | 0.43 | U | | 0.43 | U | | 0.39 | U | |
| 1,2,3-TRICHLOROBENZENE | | 0.58 | U | | 0.58 | U | | 0.53 | U | |
| 1,2,3-TRICHLOROPROPANE | | 0.9 | U | | 0.9 | U | | 0.82 | U | |
| 1,2,3-TRIMETHYLBENZENE | | 0.53 | U | | 0.53 | U | | 0.48 | U | |
| 1,2,4-TRICHLOROBENZENE | | 0.44 | U | | 0.44 | U | | 0.4 | U | |
| 1,2,4-TRIMETHYLBENZENE | | 0.67 | U | | 0.67 | U | | 0.61 | U | |
| 1,2-DIBROMO-3-CHLOROPROPANE | | 3.1 | U | | 3.1 | U | | 2.9 | U | |
| 1,2-DIBROMOETHANE | | 0.61 | U | | 0.61 | U | | 0.56 | U | |
| 1,2-DICHLOROBENZENE | | 0.39 | U | | 0.39 | U | | 0.36 | U | |
| 1,2-DICHLOROETHANE | | 0.53 | U | | 0.53 | U | | 0.48 | U | |
| 1,2-DICHLOROPROPANE | | 0.16 | U | | 0.16 | U | | 0.15 | U | |
| 1,3-DICHLOROBENZENE | | 0.8 | U | | 0.8 | U | | 0.73 | U | |
| 1,3-DICHLOROPROPANE | | 0.46 | U | | 0.46 | U | | 0.42 | U | |
| 1,4-DICHLOROBENZENE | | 0.32 | U | | 0.32 | U | | 0.29 | U | |
| 2,2-DICHLOROPROPANE | | 0.51 | UJ | C | 0.51 | UJ | C | 0.46 | UJ | C |
| 2-BUTANONE | | 4.3 | U | A | 4.4 | U | A | 3.6 | U | A |
| 2-CHLOROETHYL VINYL ETHER | | 0.74 | U | | 0.74 | U | | 0.67 | UJ | C |
| 2-CHLOROTOLUENE | | 0.35 | U | | 0.35 | U | | 0.32 | U | |
| 2-HEXANONE | | 0.78 | U | | 0.79 | U | | 0.72 | UJ | C |
| 4-CHLOROTOLUENE | | 0.3 | U | | 0.3 | U | | 0.27 | U | |
| 4-ISOPROPYLTOLUENE | | 0.47 | U | | 0.47 | U | | 0.43 | U | |
| 4-METHYL-2-PENTANONE | | 1.4 | U | | 1.4 | U | | 1.3 | UJ | C |
| ACETONE | | 6.6 | U | | 12 | J | P | 6 | UJ | C |
| BENZENE | | 1 | U | | 1 | U | | 0.95 | U | |
| BROMOBENZENE | | 0.38 | U | | 0.38 | U | | 0.35 | U | |
| BROMOCHLOROMETHANE | | 0.38 | U | | 0.38 | U | | 0.35 | U | |
| BROMODICHLOROMETHANE | | 0.33 | U | | 0.34 | U | | 0.31 | U | |
| BROMOFORM | | 0.28 | U | | 0.28 | U | | 0.25 | U | |
| BROMOMETHANE | | 0.48 | U | | 0.49 | U | | 0.44 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-10-14 | | | F-VB-UST-11-14 | | | F-VB-UST-1-14 | | | F-VB-UST-12-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-10 | | | 240-57597-11 | | | 240-57597-1 | | | 240-57597-12 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 86.6 | | | 83.5 | | | 84.3 | | | 84.1 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| CARBON DISULFIDE | | 0.66 | U | | 0.7 | U | | 0.59 | U | | 0.65 | U | |
| CARBON TETRACHLORIDE | | 0.73 | UJ | C | 0.77 | UJ | C | 0.65 | UJ | C | 0.72 | UJ | C |
| CHLOROBENZENE | | 0.6 | U | | 0.63 | U | | 0.53 | U | | 0.59 | U | |
| CHLORODIBROMOMETHANE | | 0.42 | U | | 0.44 | U | | 0.37 | U | | 0.41 | U | |
| CHLOROETHANE | | 0.49 | U | | 0.52 | U | | 0.44 | U | | 0.49 | U | |
| CHLOROFORM | | 0.4 | U | | 0.43 | U | | 0.36 | U | | 0.4 | U | |
| CHLOROMETHANE | | 0.89 | U | | 0.94 | U | | 0.8 | U | | 0.87 | U | |
| CIS-1,2-DICHLOROETHENE | | 0.31 | U | | 0.33 | U | | 0.28 | U | | 0.31 | U | |
| CIS-1,3-DICHLOROPROPENE | | 0.8 | U | | 0.85 | U | | 0.72 | U | | 0.79 | U | |
| DIBROMOMETHANE | | 0.47 | U | | 0.5 | U | | 0.42 | U | | 0.46 | U | |
| DICHLORODIFLUOROMETHANE | | 0.31 | U | | 0.33 | U | | 0.28 | U | | 0.31 | U | |
| DIISOPROPYL ETHER | | 0.66 | U | | 0.7 | U | | 0.59 | U | | 0.65 | U | |
| ETHYL TERT-BUTYL ETHER | | 0.46 | U | | 0.49 | U | | 0.41 | U | | 0.45 | U | |
| ETHYLBENZENE | | 0.36 | U | | 0.38 | U | | 0.32 | U | | 0.35 | U | |
| HEXACHLOROBUTADIENE | | 0.45 | U | | 0.48 | U | | 0.4 | U | | 0.44 | U | |
| ISOPROPYLBENZENE | | 0.22 | U | | 0.24 | U | | 0.2 | U | | 0.22 | U | |
| M+P-XYLENES | | 0.47 | U | | 0.5 | U | | 0.42 | U | | 0.46 | U | |
| METHYL TERT-BUTYL ETHER | | 0.48 | U | | 0.51 | U | | 0.43 | U | | 0.48 | U | |
| METHYLENE CHLORIDE | | 0.83 | U | | 0.88 | U | | 0.75 | U | | 0.82 | U | |
| NAPHTHALENE | | 0.39 | U | | 0.42 | U | | 0.35 | U | | 0.39 | U | |
| N-BUTYLBENZENE | | 0.66 | U | | 0.7 | U | | 0.59 | U | | 0.65 | U | |
| N-PROPYLBENZENE | | 0.67 | U | | 0.71 | U | | 0.6 | U | | 0.66 | U | |
| O-XYLENE | | 0.64 | U | | 0.68 | U | | 0.57 | U | | 0.63 | U | |
| SEC-BUTYLBENZENE | | 1 | U | | 1.1 | U | | 0.92 | U | | 1 | U | |
| STYRENE | | 0.45 | U | | 0.48 | U | | 0.4 | U | | 0.44 | U | |
| TERT-AMYL METHYL ETHER | | 0.48 | U | | 0.51 | U | | 0.43 | U | | 0.48 | U | |
| TERT-BUTYLBENZENE | | 0.75 | U | | 0.8 | U | | 0.67 | U | | 0.74 | U | |
| TERTIARY-BUTYL ALCOHOL | | 6.4 | U | | 6.8 | U | | 5.8 | U | | 6.3 | U | |
| TETRACHLOROETHENE | | 0.9 | U | | 0.95 | U | | 0.81 | U | | 0.89 | U | |
| TOLUENE | | 0.3 | U | | 0.32 | U | | 0.27 | U | | 0.3 | J | P |
| TOTAL XYLENES | | 0.61 | U | | 0.64 | U | | 0.54 | U | | 0.6 | U | |
| TRANS-1,2-DICHLOROETHENE | | 0.36 | U | | 0.38 | U | | 0.32 | U | | 0.35 | U | |
| TRANS-1,3-DICHLOROPROPENE | | 0.33 | U | | 0.35 | U | | 0.29 | U | | 0.32 | U | |
| TRICHLOROETHENE | | 0.43 | U | | 0.45 | U | | 0.38 | U | | 0.42 | U | |
| TRICHLOROFLUOROMETHANE | | 0.33 | UJ | C | 0.35 | UJ | C | 0.29 | UJ | C | 0.32 | UJ | C |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-13-14 | | | F-VB-UST-14-14 | | | F-VB-UST-2-14 | | | F-VB-UST-3-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-13 | | | 240-57597-14 | | | 240-57597-2 | | | 240-57597-3 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 85.3 | | | 85.5 | | | 85.0 | | | 82.2 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| CARBON DISULFIDE | | 0.61 | U | | 0.62 | U | | 0.65 | U | | 0.69 | U | |
| CARBON TETRACHLORIDE | | 0.67 | UJ | C | 0.68 | UJ | C | 0.71 | UJ | C | 0.76 | UJ | C |
| CHLOROBENZENE | | 0.55 | U | | 0.56 | U | | 0.58 | U | | 0.62 | U | |
| CHLORODIBROMOMETHANE | | 0.38 | U | | 0.39 | U | | 0.41 | U | | 0.43 | U | |
| CHLOROETHANE | | 0.45 | U | | 0.46 | U | | 0.48 | U | | 0.51 | U | |
| CHLOROFORM | | 0.37 | U | | 0.38 | U | | 0.39 | U | | 0.42 | U | |
| CHLOROMETHANE | | 0.82 | U | | 0.83 | U | | 0.87 | U | | 0.92 | U | |
| CIS-1,2-DICHLOROETHENE | | 0.29 | U | | 0.3 | U | | 0.31 | U | | 0.33 | U | |
| CIS-1,3-DICHLOROPROPENE | | 0.73 | U | | 0.75 | U | | 0.78 | U | | 0.83 | U | |
| DIBROMOMETHANE | | 0.43 | U | | 0.44 | U | | 0.46 | U | | 0.49 | U | |
| DICHLORODIFLUOROMETHANE | | 0.29 | U | | 0.3 | U | | 0.31 | U | | 0.33 | U | |
| DIISOPROPYL ETHER | | 0.61 | U | | 0.62 | U | | 0.65 | U | | 0.69 | U | |
| ETHYL TERT-BUTYL ETHER | | 0.42 | U | | 0.43 | U | | 0.45 | U | | 0.48 | U | |
| ETHYLBENZENE | | 0.33 | U | | 0.34 | U | | 0.35 | U | | 0.37 | U | |
| HEXACHLOROBUTADIENE | | 0.41 | U | | 0.42 | U | | 0.44 | U | | 0.46 | U | |
| ISOPROPYLBENZENE | | 0.21 | U | | 0.21 | U | | 0.22 | U | | 0.23 | U | |
| M+P-XYLENES | | 0.43 | U | | 0.44 | U | | 0.46 | U | | 0.49 | U | |
| METHYL TERT-BUTYL ETHER | | 0.44 | U | | 0.45 | U | | 0.47 | U | | 0.5 | U | |
| METHYLENE CHLORIDE | | 0.76 | U | | 0.78 | U | | 0.81 | U | | 0.86 | U | |
| NAPHTHALENE | | 0.36 | U | | 0.37 | U | | 0.38 | U | | 0.41 | U | |
| N-BUTYLBENZENE | | 0.61 | U | | 0.62 | U | | 0.65 | U | | 0.69 | U | |
| N-PROPYLBENZENE | | 0.62 | U | | 0.63 | U | | 0.66 | U | | 0.7 | U | |
| O-XYLENE | | 0.59 | U | | 0.6 | U | | 0.63 | U | | 0.66 | U | |
| SEC-BUTYLBENZENE | | 0.94 | U | | 0.96 | U | | 1 | U | | 1.1 | U | |
| STYRENE | | 0.41 | U | | 0.42 | U | | 0.44 | U | | 0.46 | U | |
| TERT-AMYL METHYL ETHER | | 0.44 | U | | 0.45 | U | | 0.47 | U | | 0.5 | U | |
| TERT-BUTYLBENZENE | | 0.69 | U | | 0.71 | U | | 0.74 | U | | 0.78 | U | |
| TERTIARY-BUTYL ALCOHOL | | 5.9 | U | | 6 | U | | 6.3 | U | | 6.6 | U | |
| TETRACHLOROETHENE | | 0.83 | U | | 0.84 | U | | 0.88 | U | | 0.93 | U | |
| TOLUENE | | 0.28 | U | | 0.29 | J | P | 0.3 | U | | 0.31 | U | |
| TOTAL XYLENES | | 0.56 | U | | 0.57 | U | | 0.59 | U | | 0.63 | U | |
| TRANS-1,2-DICHLOROETHENE | | 0.33 | U | | 0.34 | U | | 0.35 | U | | 0.37 | U | |
| TRANS-1,3-DICHLOROPROPENE | | 0.3 | U | | 0.31 | U | | 0.32 | U | | 0.34 | U | |
| TRICHLOROETHENE | | 0.39 | U | | 0.4 | U | | 0.42 | U | | 0.44 | U | |
| TRICHLOROFLUOROMETHANE | | 0.3 | UJ | C | 0.31 | UJ | C | 0.32 | UJ | C | 0.34 | UJ | C |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-4-14 | | | F-VB-UST-5-14 | | | F-VB-UST-6-14 | | | F-VB-UST-7-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-4 | | | 240-57597-5 | | | 240-57597-6 | | | 240-57597-7 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 84.4 | | | 84.8 | | | 83.5 | | | 83.8 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| CARBON DISULFIDE | | 0.64 | U | | 0.68 | U | | 0.68 | U | | 0.66 | U | |
| CARBON TETRACHLORIDE | | 0.71 | UJ | C | 0.75 | UJ | C | 0.75 | UJ | C | 0.73 | UJ | C |
| CHLOROBENZENE | | 0.58 | U | | 0.61 | U | | 0.61 | U | | 0.6 | U | |
| CHLORODIBROMOMETHANE | | 0.4 | U | | 0.43 | U | | 0.43 | U | | 0.42 | U | |
| CHLOROETHANE | | 0.48 | U | | 0.51 | U | | 0.51 | U | | 0.49 | U | |
| CHLOROFORM | | 0.39 | U | | 0.42 | U | | 0.42 | U | | 0.4 | U | |
| CHLOROMETHANE | | 0.86 | U | | 0.92 | U | | 0.91 | U | | 0.89 | U | |
| CIS-1,2-DICHLOROETHENE | | 0.31 | U | | 0.32 | U | | 0.32 | U | | 0.31 | U | |
| CIS-1,3-DICHLOROPROPENE | | 0.77 | U | | 0.82 | U | | 0.82 | U | | 0.8 | U | |
| DIBROMOMETHANE | | 0.46 | U | | 0.49 | U | | 0.49 | U | | 0.47 | U | |
| DICHLORODIFLUOROMETHANE | | 0.31 | U | | 0.32 | U | | 0.32 | U | | 0.31 | U | |
| DIISOPROPYL ETHER | | 0.64 | U | | 0.68 | U | | 0.68 | U | | 0.66 | U | |
| ETHYL TERT-BUTYL ETHER | | 0.45 | U | | 0.47 | U | | 0.47 | U | | 0.46 | U | |
| ETHYLBENZENE | | 0.35 | U | | 0.37 | U | | 0.37 | U | | 0.36 | U | |
| HEXACHLOROBUTADIENE | | 0.44 | U | | 0.46 | U | | 0.46 | U | | 0.45 | U | |
| ISOPROPYLBENZENE | | 0.22 | U | | 0.23 | U | | 0.23 | U | | 0.22 | U | |
| M+P-XYLENES | | 0.46 | U | | 0.49 | U | | 0.49 | U | | 0.47 | U | |
| METHYL TERT-BUTYL ETHER | | 0.47 | U | | 0.5 | U | | 0.5 | U | | 0.48 | U | |
| METHYLENE CHLORIDE | | 0.81 | U | | 0.86 | U | | 0.86 | U | | 0.83 | U | |
| NAPHTHALENE | | 0.38 | U | | 0.41 | J | P | 0.45 | J | P | 0.39 | U | |
| N-BUTYLBENZENE | | 0.64 | U | | 0.68 | U | | 0.68 | U | | 0.66 | U | |
| N-PROPYLBENZENE | | 0.65 | U | | 0.7 | U | | 0.69 | U | | 0.67 | U | |
| O-XYLENE | | 0.62 | U | | 0.66 | U | | 0.66 | U | | 0.64 | U | |
| SEC-BUTYLBENZENE | | 0.99 | U | | 1.1 | U | | 1.1 | U | | 1 | U | |
| STYRENE | | 0.44 | U | | 0.46 | U | | 0.46 | U | | 0.45 | U | |
| TERT-AMYL METHYL ETHER | | 0.47 | U | | 0.5 | U | | 0.5 | U | | 0.48 | U | |
| TERT-BUTYLBENZENE | | 0.73 | U | | 0.78 | U | | 0.77 | U | | 0.75 | U | |
| TERTIARY-BUTYL ALCOHOL | | 6.2 | U | | 6.6 | U | | 6.6 | U | | 6.4 | U | |
| TETRACHLOROETHENE | | 0.87 | U | | 0.93 | U | | 0.93 | U | | 0.9 | U | |
| TOLUENE | | 0.29 | U | | 0.31 | U | | 0.31 | U | | 0.3 | U | |
| TOTAL XYLENES | | 0.59 | U | | 0.63 | U | | 0.62 | U | | 0.61 | U | |
| TRANS-1,2-DICHLOROETHENE | | 0.35 | U | | 0.37 | U | | 0.37 | U | | 0.36 | U | |
| TRANS-1,3-DICHLOROPROPENE | | 0.32 | U | | 0.34 | U | | 0.34 | U | | 0.33 | U | |
| TRICHLOROETHENE | | 0.41 | U | | 0.44 | U | | 0.44 | U | | 0.43 | U | |
| TRICHLOROFLUOROMETHANE | | 0.32 | UJ | C | 0.34 | UJ | C | 0.34 | UJ | C | 0.33 | UJ | C |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-8-14 | | | F-VB-UST-9-14 | | | F-VB-UST-DUP | | |
|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-8 | | | 240-57597-9 | | | 240-57597-15 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 82.0 | | | 80.9 | | | 85.5 | | |
| | DUP_OF | | | | | | | F-VB-UST-2-14 | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| CARBON DISULFIDE | | 0.68 | U | | 0.68 | U | | 0.62 | U | |
| CARBON TETRACHLORIDE | | 0.75 | UJ | C | 0.75 | UJ | C | 0.68 | U | |
| CHLOROBENZENE | | 0.61 | U | | 0.61 | U | | 0.56 | U | |
| CHLORODIBROMOMETHANE | | 0.43 | U | | 0.43 | U | | 0.39 | U | |
| CHLOROETHANE | | 0.51 | U | | 0.51 | U | | 0.46 | U | |
| CHLOROFORM | | 0.41 | U | | 0.42 | U | | 0.38 | U | |
| CHLOROMETHANE | | 0.91 | U | | 0.91 | U | | 0.83 | U | |
| CIS-1,2-DICHLOROETHENE | | 0.32 | U | | 0.32 | U | | 0.29 | U | |
| CIS-1,3-DICHLOROPROPENE | | 0.82 | U | | 0.82 | U | | 0.75 | U | |
| DIBROMOMETHANE | | 0.48 | U | | 0.49 | U | | 0.44 | U | |
| DICHLORODIFLUOROMETHANE | | 0.32 | U | | 0.32 | U | | 0.29 | U | |
| DIISOPROPYL ETHER | | 0.68 | U | | 0.68 | U | | 0.62 | U | |
| ETHYL TERT-BUTYL ETHER | | 0.47 | U | | 0.47 | U | | 0.43 | U | |
| ETHYLBENZENE | | 0.37 | U | | 0.37 | U | | 0.34 | U | |
| HEXACHLOROBUTADIENE | | 0.46 | U | | 0.46 | U | | 0.42 | U | |
| ISOPROPYLBENZENE | | 0.23 | U | | 0.23 | U | | 0.21 | U | |
| M+P-XYLENES | | 0.48 | U | | 0.49 | U | | 0.44 | U | |
| METHYL TERT-BUTYL ETHER | | 0.5 | U | | 0.5 | U | | 0.45 | U | |
| METHYLENE CHLORIDE | | 0.85 | U | | 0.85 | U | | 0.78 | U | |
| NAPHTHALENE | | 0.4 | U | | 0.4 | U | | 0.37 | U | |
| N-BUTYLBENZENE | | 0.68 | U | | 0.68 | U | | 0.62 | U | |
| N-PROPYLBENZENE | | 0.69 | U | | 0.69 | U | | 0.63 | U | |
| O-XYLENE | | 0.66 | U | | 0.66 | U | | 0.6 | U | |
| SEC-BUTYLBENZENE | | 1 | U | | 1.1 | U | | 0.96 | U | |
| STYRENE | | 0.46 | U | | 0.46 | U | | 0.42 | U | |
| TERT-AMYL METHYL ETHER | | 0.5 | U | | 0.5 | U | | 0.45 | U | |
| TERT-BUTYLBENZENE | | 0.77 | U | | 0.77 | U | | 0.71 | UJ | C |
| TERTIARY-BUTYL ALCOHOL | | 6.6 | U | | 6.6 | U | | 6 | U | |
| TETRACHLOROETHENE | | 0.92 | U | | 0.92 | U | | 0.84 | U | |
| TOLUENE | | 0.33 | J | P | 0.36 | J | P | 0.28 | U | |
| TOTAL XYLENES | | 0.62 | U | | 0.62 | U | | 0.57 | U | |
| TRANS-1,2-DICHLOROETHENE | | 0.37 | U | | 0.37 | U | | 0.34 | U | |
| TRANS-1,3-DICHLOROPROPENE | | 0.33 | U | | 0.34 | U | | 0.31 | U | |
| TRICHLOROETHENE | | 0.44 | U | | 0.44 | U | | 0.4 | U | |
| TRICHLOROFLUOROMETHANE | | 0.33 | UJ | C | 0.34 | UJ | C | 0.31 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-10-14 | | | F-VB-UST-11-14 | | | F-VB-UST-1-14 | | | F-VB-UST-12-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-10 | | | 240-57597-11 | | | 240-57597-1 | | | 240-57597-12 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 86.6 | | | 83.5 | | | 84.3 | | | 84.1 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| VINYL ACETATE | | 0.74 | U | | 0.79 | U | | 0.66 | U | | 0.73 | U | |
| VINYL CHLORIDE | | 0.34 | U | | 0.36 | U | | 0.3 | U | | 0.33 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-13-14 | | | F-VB-UST-14-14 | | | F-VB-UST-2-14 | | | F-VB-UST-3-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-13 | | | 240-57597-14 | | | 240-57597-2 | | | 240-57597-3 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 85.3 | | | 85.5 | | | 85.0 | | | 82.2 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| VINYL ACETATE | | 0.68 | U | | 0.7 | U | | 0.72 | U | | 0.77 | U | |
| VINYL CHLORIDE | | 0.31 | U | | 0.32 | U | | 0.33 | U | | 0.35 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-4-14 | | | F-VB-UST-5-14 | | | F-VB-UST-6-14 | | | F-VB-UST-7-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-4 | | | 240-57597-5 | | | 240-57597-6 | | | 240-57597-7 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 84.4 | | | 84.8 | | | 83.5 | | | 83.8 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| VINYL ACETATE | | 0.72 | U | | 0.76 | U | | 0.76 | U | | 0.74 | U | |
| VINYL CHLORIDE | | 0.33 | U | | 0.35 | U | | 0.35 | U | | 0.34 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-8-14 | | | F-VB-UST-9-14 | | | F-VB-UST-DUP | | |
|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-8 | | | 240-57597-9 | | | 240-57597-15 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 82.0 | | | 80.9 | | | 85.5 | | |
| | DUP_OF | | | | | | | F-VB-UST-2-14 | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| VINYL ACETATE | | 0.76 | U | | 0.76 | U | | 0.7 | U | |
| VINYL CHLORIDE | | 0.35 | U | | 0.35 | U | | 0.32 | U | |

| | | | | |
|---|---|---|---|---|
| **PROJ_NO: 07336** | NSAMPLE | TB-110515 | | |
| **SDG: 240-57597-1** | LAB_ID | 240-57597-16 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | |
| **MEDIA: WATER** | QC_TYPE | NM | | |
| | UNITS | UG/L | | |
| | PCT_SOLIDS | 0.0 | | |
| | DUP_OF | | | |
| PARAMETER | | RESULT | VQL | QLCD |
| 1,1,1,2-TETRACHLOROETHANE | | 0.28 | U | |
| 1,1,1-TRICHLOROETHANE | | 0.44 | U | |
| 1,1,2,2-TETRACHLOROETHANE | | 0.22 | U | |
| 1,1,2-TRICHLOROTRIFLUOROETHANE | | 0.45 | U | |
| 1,1-DICHLOROETHANE | | 0.3 | U | |
| 1,1-DICHLOROETHENE | | 0.45 | U | |
| 1,1-DICHLOROPROPENE | | 0.42 | U | |
| 1,2,3-TRICHLOROBENZENE | | 0.37 | U | |
| 1,2,3-TRICHLOROPROPANE | | 0.44 | U | |
| 1,2,3-TRIMETHYLBENZENE | | 0.47 | U | |
| 1,2,4-TRICHLOROBENZENE | | 0.32 | U | |
| 1,2,4-TRIMETHYLBENZENE | | 0.41 | U | |
| 1,2-DIBROMO-3-CHLOROPROPANE | | 0.82 | U | |
| 1,2-DIBROMOETHANE | | 0.32 | U | |
| 1,2-DICHLOROBENZENE | | 0.25 | U | |
| 1,2-DICHLOROETHANE | | 0.23 | U | |
| 1,2-DICHLOROPROPANE | | 0.25 | U | |
| 1,3-DICHLOROBENZENE | | 0.19 | U | |
| 1,3-DICHLOROPROPANE | | 0.19 | U | |
| 1,4-DICHLOROBENZENE | | 0.27 | U | |
| 2,2-DICHLOROPROPANE | | 0.26 | U | |
| 2-BUTANONE | | 0.53 | U | |
| 2-CHLOROETHYL VINYL ETHER | | 0.59 | U | |
| 2-CHLOROTOLUENE | | 0.4 | U | |
| 2-HEXANONE | | 0.48 | U | |
| 4-CHLOROTOLUENE | | 0.29 | U | |
| 4-ISOPROPYLTOLUENE | | 0.43 | U | |
| 4-METHYL-2-PENTANONE | | 0.99 | U | |
| ACETONE | | 0.94 | U | |
| BENZENE | | 0.35 | U | |
| BROMOBENZENE | | 0.35 | U | |
| BROMOCHLOROMETHANE | | 0.5 | U | |
| BROMODICHLOROMETHANE | | 0.29 | U | |
| BROMOFORM | | 0.56 | U | |
| BROMOMETHANE | | 0.44 | U | |

| | | | | |
|---|---|---|---|---|
| **PROJ_NO: 07336** | NSAMPLE | TB-110515 | | |
| **SDG: 240-57597-1** | LAB_ID | 240-57597-16 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | |
| **MEDIA: WATER** | QC_TYPE | NM | | |
| | UNITS | UG/L | | |
| | PCT_SOLIDS | 0.0 | | |
| | DUP_OF | | | |
| PARAMETER | | RESULT | VQL | QLCD |
| CARBON DISULFIDE | | 0.38 | U | |
| CARBON TETRACHLORIDE | | 0.43 | U | |
| CHLOROBENZENE | | 0.25 | U | |
| CHLORODIBROMOMETHANE | | 0.43 | U | |
| CHLOROETHANE | | 0.32 | U | |
| CHLOROFORM | | 0.25 | U | |
| CHLOROMETHANE | | 0.44 | U | |
| CIS-1,2-DICHLOROETHENE | | 0.26 | U | |
| CIS-1,3-DICHLOROPROPENE | | 0.46 | U | |
| DIBROMOMETHANE | | 0.42 | U | |
| DICHLORODIFLUOROMETHANE | | 0.32 | U | |
| DIISOPROPYL ETHER | | 0.5 | U | |
| ETHYL TERT-BUTYL ETHER | | 0.23 | U | |
| ETHYLBENZENE | | 0.25 | U | |
| HEXACHLOROBUTADIENE | | 0.35 | UJ | C |
| ISOPROPYLBENZENE | | 0.35 | U | |
| M+P-XYLENES | | 0.24 | U | |
| METHYL TERT-BUTYL ETHER | | 0.2 | U | |
| METHYLENE CHLORIDE | | 0.33 | U | |
| NAPHTHALENE | | 0.45 | U | |
| N-BUTYLBENZENE | | 0.31 | U | |
| N-PROPYLBENZENE | | 0.4 | U | |
| O-XYLENE | | 0.25 | U | |
| SEC-BUTYLBENZENE | | 0.48 | U | |
| STYRENE | | 0.45 | U | |
| TERT-AMYL METHYL ETHER | | 0.3 | U | |
| TERT-BUTYLBENZENE | | 0.41 | U | |
| TERTIARY-BUTYL ALCOHOL | | 4.9 | U | |
| TETRACHLOROETHENE | | 0.31 | U | |
| TOLUENE | | 0.23 | U | |
| TOTAL XYLENES | | 0.52 | U | |
| TRANS-1,2-DICHLOROETHENE | | 0.3 | U | |
| TRANS-1,3-DICHLOROPROPENE | | 0.56 | U | |
| TRICHLOROETHENE | | 0.22 | U | |
| TRICHLOROFLUOROMETHANE | | 0.49 | U | |

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|---|---|---|---|---|
| **PROJ_NO: 07336** | NSAMPLE | TB-110515 | | |
| **SDG: 240-57597-1** | LAB_ID | 240-57597-16 | | |
| **FRACTION: OV** | SAMP_DATE | 11/5/2015 | | |
| **MEDIA: WATER** | QC_TYPE | NM | | |
| | UNITS | UG/L | | |
| | PCT_SOLIDS | 0.0 | | |
| | DUP_OF | | | |
| PARAMETER | | RESULT | VQL | QLCD |
| VINYL ACETATE | | 0.41 | U | |
| VINYL CHLORIDE | | 0.29 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-10-14 | | | F-VB-UST-11-14 | | | F-VB-UST-1-14 | | | F-VB-UST-12-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-10 | | | 240-57597-11 | | | 240-57597-1 | | | 240-57597-12 | | |
| **FRACTION: OS** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 86.6 | | | 83.5 | | | 84.3 | | | 84.1 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| 1,1-BIPHENYL | | 4.1 | U | | 4.2 | U | | 4.2 | U | | 4.2 | U | |
| 2,2'-OXYBIS(1-CHLOROPROPANE) | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| 2,4,5-TRICHLOROPHENOL | | 29 | U | | 30 | U | | 30 | U | | 30 | U | |
| 2,4,6-TRICHLOROPHENOL | | 10 | U | | 11 | U | | 11 | U | | 11 | U | |
| 2,4-DICHLOROPHENOL | | 23 | U | | 24 | U | | 24 | U | | 24 | U | |
| 2,4-DIMETHYLPHENOL | | 23 | U | | 24 | U | | 24 | U | | 24 | U | |
| 2,4-DINITROPHENOL | | 24 | U | | 25 | U | | 25 | U | | 25 | U | |
| 2,4-DINITROTOLUENE | | 20 | U | | 20 | U | | 20 | U | | 20 | U | |
| 2,6-DINITROTOLUENE | | 24 | U | | 25 | U | | 25 | U | | 25 | U | |
| 2-CHLORONAPHTHALENE | | 0.52 | U | | 0.54 | U | | 0.53 | U | | 0.54 | U | |
| 2-CHLOROPHENOL | | 9.5 | U | | 9.9 | U | | 9.7 | U | | 9.8 | U | |
| 2-METHYLNAPHTHALENE | | 0.58 | U | | 0.6 | U | | 0.59 | U | | 0.6 | U | |
| 2-METHYLPHENOL | | 13 | U | | 13 | U | | 13 | U | | 13 | U | |
| 2-NITROANILINE | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| 2-NITROPHENOL | | 9.7 | U | | 10 | U | | 9.9 | U | | 9.9 | U | |
| 3&4-METHYLPHENOL | | 23 | U | | 24 | U | | 24 | U | | 24 | U | |
| 3,3'-DICHLOROBENZIDINE | | 21 | U | | 22 | U | | 21 | U | | 22 | U | |
| 3-NITROANILINE | | 19 | U | | 19 | U | | 19 | U | | 19 | U | |
| 4,6-DINITRO-2-METHYLPHENOL | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| 4-BROMOPHENYL PHENYL ETHER | | 15 | U | | 16 | U | | 15 | U | | 16 | U | |
| 4-CHLORO-3-METHYLPHENOL | | 24 | U | | 25 | U | | 25 | U | | 25 | U | |
| 4-CHLOROANILINE | | 20 | U | | 20 | U | | 20 | U | | 20 | U | |
| 4-CHLOROPHENYL PHENYL ETHER | | 15 | U | | 16 | U | | 15 | U | | 16 | U | |
| 4-NITROANILINE | | 30 | U | | 31 | U | | 31 | U | | 31 | U | |
| 4-NITROPHENOL | | 20 | U | | 20 | U | | 20 | U | | 20 | U | |
| ACENAPHTHENE | | 0.88 | U | | 0.91 | U | | 0.9 | U | | 0.91 | U | |
| ACENAPHTHYLENE | | 0.41 | U | | 0.42 | U | | 0.42 | U | | 0.42 | U | |
| ACETOPHENONE | | 11 | U | | 38 | J | P | 16 | J | P | 11 | U | |
| ANTHRACENE | | 0.91 | U | | 0.94 | U | | 0.93 | U | | 0.93 | U | |
| ATRAZINE | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| BENZALDEHYDE | | 14 | UJ | C | 19 | J | CP | 14 | UJ | C | 14 | UJ | C |
| BENZO(A)ANTHRACENE | | 0.73 | U | | 0.76 | U | | 0.75 | U | | 0.75 | U | |
| BENZO(A)PYRENE | | 0.74 | U | | 0.77 | U | | 0.76 | U | | 0.77 | U | |
| BENZO(B)FLUORANTHENE | | 0.69 | U | | 0.71 | U | | 0.7 | U | | 0.71 | U | |
| BENZO(G,H,I)PERYLENE | | 0.41 | U | | 0.42 | U | | 0.42 | U | | 0.42 | U | |

| PROJ_NO: 07336 | NSAMPLE | F-VB-UST-13-14 | | | F-VB-UST-14-14 | | | F-VB-UST-2-14 | | | F-VB-UST-3-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| SDG: 240-57597-1 | LAB_ID | 240-57597-13 | | | 240-57597-14 | | | 240-57597-2 | | | 240-57597-3 | | |
| FRACTION: OS | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| MEDIA: SOIL | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 85.3 | | | 85.5 | | | 85.0 | | | 82.2 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| 1,1-BIPHENYL | | 4.1 | U | | 4.1 | U | | 4.1 | U | | 4.3 | U | |
| 2,2'-OXYBIS(1-CHLOROPROPANE) | | 11 | U | | 11 | U | | 11 | U | | 12 | U | |
| 2,4,5-TRICHLOROPHENOL | | 29 | U | | 30 | U | | 29 | U | | 30 | U | |
| 2,4,6-TRICHLOROPHENOL | | 10 | U | | 11 | U | | 10 | U | | 11 | U | |
| 2,4-DICHLOROPHENOL | | 23 | U | | 24 | U | | 23 | U | | 24 | U | |
| 2,4-DIMETHYLPHENOL | | 23 | U | | 24 | U | | 23 | U | | 24 | U | |
| 2,4-DINITROPHENOL | | 24 | U | | 25 | U | | 25 | U | | 26 | U | |
| 2,4-DINITROTOLUENE | | 20 | U | | 20 | U | | 20 | U | | 21 | U | |
| 2,6-DINITROTOLUENE | | 24 | U | | 25 | U | | 25 | U | | 26 | U | |
| 2-CHLORONAPHTHALENE | | 0.52 | U | | 0.53 | U | | 0.53 | U | | 0.55 | U | |
| 2-CHLOROPHENOL | | 9.5 | U | | 9.7 | U | | 9.6 | U | | 10 | U | |
| 2-METHYLNAPHTHALENE | | 0.58 | U | | 0.59 | U | | 0.58 | U | | 0.61 | U | |
| 2-METHYLPHENOL | | 13 | U | | 13 | U | | 13 | U | | 13 | U | |
| 2-NITROANILINE | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| 2-NITROPHENOL | | 9.6 | U | | 9.8 | U | | 9.7 | U | | 10 | U | |
| 3&4-METHYLPHENOL | | 23 | U | | 24 | U | | 23 | U | | 24 | U | |
| 3,3'-DICHLOROBENZIDINE | | 21 | U | | 21 | U | | 21 | U | | 22 | U | |
| 3-NITROANILINE | | 19 | U | | 19 | U | | 19 | U | | 19 | U | |
| 4,6-DINITRO-2-METHYLPHENOL | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| 4-BROMOPHENYL PHENYL ETHER | | 15 | U | | 15 | U | | 15 | U | | 16 | U | |
| 4-CHLORO-3-METHYLPHENOL | | 24 | U | | 25 | U | | 25 | U | | 26 | U | |
| 4-CHLOROANILINE | | 20 | U | | 20 | U | | 20 | U | | 21 | U | |
| 4-CHLOROPHENYL PHENYL ETHER | | 15 | U | | 15 | U | | 15 | U | | 16 | U | |
| 4-NITROANILINE | | 30 | U | | 31 | U | | 30 | U | | 32 | U | |
| 4-NITROPHENOL | | 20 | U | | 20 | U | | 20 | U | | 21 | U | |
| ACENAPHTHENE | | 0.88 | U | | 0.9 | U | | 0.89 | U | | 0.92 | U | |
| ACENAPHTHYLENE | | 0.41 | U | | 0.41 | U | | 0.41 | U | | 0.43 | U | |
| ACETOPHENONE | | 11 | U | | 11 | J | P | 11 | U | | 11 | U | |
| ANTHRACENE | | 0.9 | U | | 0.92 | U | | 0.91 | U | | 0.95 | U | |
| ATRAZINE | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| BENZALDEHYDE | | 14 | UJ | C | 14 | UJ | C | 14 | UJ | C | 15 | UJ | C |
| BENZO(A)ANTHRACENE | | 0.73 | U | | 0.75 | U | | 0.74 | U | | 0.77 | U | |
| BENZO(A)PYRENE | | 0.74 | U | | 0.76 | U | | 0.75 | U | | 0.78 | U | |
| BENZO(B)FLUORANTHENE | | 0.68 | U | | 0.7 | U | | 0.69 | U | | 0.72 | U | |
| BENZO(G,H,I)PERYLENE | | 0.41 | U | | 0.41 | U | | 0.41 | U | | 0.43 | U | |

| PROJ_NO: 07336 | NSAMPLE | F-VB-UST-4-14 | | | F-VB-UST-5-14 | | | F-VB-UST-6-14 | | | F-VB-UST-7-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| SDG: 240-57597-1 | LAB_ID | 240-57597-4 | | | 240-57597-5 | | | 240-57597-6 | | | 240-57597-7 | | |
| FRACTION: OS | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| MEDIA: SOIL | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 84.4 | | | 84.8 | | | 83.5 | | | 83.8 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| 1,1-BIPHENYL | | 4.2 | U | | 4.1 | U | | 4.2 | U | | 4.2 | U | |
| 2,2'-OXYBIS(1-CHLOROPROPANE) | | 11 | U | | 11 | U | | 11 | U | | 12 | U | |
| 2,4,5-TRICHLOROPHENOL | | 30 | U | | 29 | U | | 30 | U | | 30 | U | |
| 2,4,6-TRICHLOROPHENOL | | 11 | U | | 10 | U | | 11 | U | | 11 | U | |
| 2,4-DICHLOROPHENOL | | 24 | U | | 23 | U | | 24 | U | | 24 | U | |
| 2,4-DIMETHYLPHENOL | | 24 | U | | 23 | U | | 24 | U | | 24 | U | |
| 2,4-DINITROPHENOL | | 25 | U | | 25 | U | | 25 | U | | 25 | U | |
| 2,4-DINITROTOLUENE | | 20 | U | | 20 | U | | 20 | U | | 21 | U | |
| 2,6-DINITROTOLUENE | | 25 | U | | 25 | U | | 25 | U | | 25 | U | |
| 2-CHLORONAPHTHALENE | | 0.54 | U | | 0.53 | U | | 0.54 | U | | 0.55 | U | |
| 2-CHLOROPHENOL | | 9.8 | U | | 9.6 | U | | 9.8 | U | | 9.9 | U | |
| 2-METHYLNAPHTHALENE | | 0.6 | U | | 0.58 | U | | 0.6 | U | | 0.61 | U | |
| 2-METHYLPHENOL | | 13 | U | | 13 | U | | 13 | U | | 13 | U | |
| 2-NITROANILINE | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| 2-NITROPHENOL | | 9.9 | U | | 9.7 | U | | 9.9 | U | | 10 | U | |
| 3&4-METHYLPHENOL | | 24 | U | | 23 | U | | 24 | U | | 24 | U | |
| 3,3'-DICHLOROBENZIDINE | | 21 | U | | 21 | U | | 21 | U | | 22 | U | |
| 3-NITROANILINE | | 19 | U | | 19 | U | | 19 | U | | 19 | U | |
| 4,6-DINITRO-2-METHYLPHENOL | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| 4-BROMOPHENYL PHENYL ETHER | | 15 | U | | 15 | U | | 15 | U | | 16 | U | |
| 4-CHLORO-3-METHYLPHENOL | | 25 | U | | 25 | U | | 25 | U | | 25 | U | |
| 4-CHLOROANILINE | | 20 | U | | 20 | U | | 20 | U | | 21 | U | |
| 4-CHLOROPHENYL PHENYL ETHER | | 15 | U | | 15 | U | | 15 | U | | 16 | U | |
| 4-NITROANILINE | | 31 | U | | 30 | U | | 31 | U | | 32 | U | |
| 4-NITROPHENOL | | 20 | U | | 20 | U | | 20 | U | | 21 | U | |
| ACENAPHTHENE | | 0.91 | U | | 0.89 | U | | 0.91 | U | | 0.92 | U | |
| ACENAPHTHYLENE | | 0.42 | U | | 0.41 | U | | 0.42 | U | | 0.42 | U | |
| ACETOPHENONE | | 13 | J | P | 11 | U | | 11 | U | | 11 | U | |
| ANTHRACENE | | 0.93 | U | | 0.91 | U | | 0.93 | U | | 0.95 | U | |
| ATRAZINE | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| BENZALDEHYDE | | 14 | UJ | C | 14 | UJ | C | 14 | UJ | C | 15 | UJ | C |
| BENZO(A)ANTHRACENE | | 0.75 | U | | 0.74 | U | | 0.75 | U | | 0.76 | U | |
| BENZO(A)PYRENE | | 0.76 | U | | 0.75 | U | | 0.76 | U | | 0.78 | U | |
| BENZO(B)FLUORANTHENE | | 0.7 | U | | 0.69 | U | | 0.7 | U | | 0.72 | U | |
| BENZO(G,H,I)PERYLENE | | 0.42 | U | | 0.41 | U | | 0.42 | U | | 0.42 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-8-14 | | | F-VB-UST-9-14 | | | F-VB-UST-DUP | | |
|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-8 | | | 240-57597-9 | | | 240-57597-15 | | |
| **FRACTION: OS** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 82.0 | | | 80.9 | | | 85.5 | | |
| | DUP_OF | | | | | | | F-VB-UST-2-14 | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| 1,1-BIPHENYL | | 4.2 | U | | 4.4 | U | | 4.1 | U | |
| 2,2'-OXYBIS(1-CHLOROPROPANE) | | 12 | U | | 12 | U | | 11 | U | |
| 2,4,5-TRICHLOROPHENOL | | 30 | U | | 31 | U | | 29 | U | |
| 2,4,6-TRICHLOROPHENOL | | 11 | U | | 11 | U | | 10 | U | |
| 2,4-DICHLOROPHENOL | | 24 | U | | 25 | U | | 23 | U | |
| 2,4-DIMETHYLPHENOL | | 24 | U | | 25 | U | | 23 | U | |
| 2,4-DINITROPHENOL | | 25 | U | | 26 | U | | 24 | U | |
| 2,4-DINITROTOLUENE | | 21 | U | | 21 | U | | 20 | U | |
| 2,6-DINITROTOLUENE | | 25 | U | | 26 | U | | 24 | U | |
| 2-CHLORONAPHTHALENE | | 0.55 | U | | 0.56 | U | | 0.52 | U | |
| 2-CHLOROPHENOL | | 9.9 | U | | 10 | U | | 9.6 | U | |
| 2-METHYLNAPHTHALENE | | 0.61 | U | | 0.62 | U | | 0.58 | U | |
| 2-METHYLPHENOL | | 13 | U | | 14 | U | | 13 | U | |
| 2-NITROANILINE | | 11 | U | | 11 | U | | 11 | U | |
| 2-NITROPHENOL | | 10 | U | | 10 | U | | 9.7 | U | |
| 3&4-METHYLPHENOL | | 24 | U | | 25 | U | | 23 | U | |
| 3,3'-DICHLOROBENZIDINE | | 22 | U | | 22 | U | | 21 | U | |
| 3-NITROANILINE | | 19 | U | | 20 | U | | 19 | U | |
| 4,6-DINITRO-2-METHYLPHENOL | | 11 | U | | 11 | U | | 11 | U | |
| 4-BROMOPHENYL PHENYL ETHER | | 16 | U | | 16 | U | | 15 | U | |
| 4-CHLORO-3-METHYLPHENOL | | 25 | U | | 26 | U | | 24 | U | |
| 4-CHLOROANILINE | | 21 | U | | 21 | U | | 20 | U | |
| 4-CHLOROPHENYL PHENYL ETHER | | 16 | U | | 16 | U | | 15 | U | |
| 4-NITROANILINE | | 31 | U | | 32 | U | | 30 | U | |
| 4-NITROPHENOL | | 21 | U | | 21 | U | | 20 | U | |
| ACENAPHTHENE | | 0.92 | U | | 0.95 | U | | 0.89 | U | |
| ACENAPHTHYLENE | | 0.42 | U | | 0.44 | U | | 0.41 | U | |
| ACETOPHENONE | | 11 | U | | 13 | J | P | 11 | U | |
| ANTHRACENE | | 0.94 | U | | 0.97 | U | | 0.91 | U | |
| ATRAZINE | | 11 | U | | 11 | U | | 11 | U | |
| BENZALDEHYDE | | 15 | UJ | C | 15 | UJ | C | 14 | UJ | C |
| BENZO(A)ANTHRACENE | | 0.76 | U | | 0.79 | U | | 0.73 | U | |
| BENZO(A)PYRENE | | 0.78 | U | | 0.8 | U | | 0.75 | U | |
| BENZO(B)FLUORANTHENE | | 0.71 | U | | 0.74 | U | | 0.69 | U | |
| BENZO(G,H,I)PERYLENE | | 0.42 | U | | 0.44 | U | | 0.41 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-10-14 | | | F-VB-UST-11-14 | | | F-VB-UST-1-14 | | | F-VB-UST-12-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-10 | | | 240-57597-11 | | | 240-57597-1 | | | 240-57597-12 | | |
| **FRACTION: OS** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 86.6 | | | 83.5 | | | 84.3 | | | 84.1 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| BENZO(K)FLUORANTHENE | | 0.79 | U | | 0.82 | U | | 0.81 | U | | 0.81 | U | |
| BIS(2-CHLOROETHOXY)METHANE | | 26 | U | | 26 | U | | 26 | U | | 26 | U | |
| BIS(2-CHLOROETHYL)ETHER | | 2.3 | U | | 2.4 | U | | 2.4 | U | | 2.4 | U | |
| BIS(2-ETHYLHEXYL)PHTHALATE | | 29 | J | P | 23 | U | | 31 | J | P | 23 | U | |
| BUTYL BENZYL PHTHALATE | | 12 | U | | 12 | U | | 12 | U | | 12 | U | |
| CAPROLACTAM | | 43 | U | | 45 | U | | 44 | U | | 44 | U | |
| CARBAZOLE | | 31 | U | | 32 | U | | 32 | U | | 32 | U | |
| CHRYSENE | | 1.3 | U | | 1.3 | U | | 1.3 | U | | 1.3 | U | |
| DIBENZO(A,H)ANTHRACENE | | 0.77 | U | | 0.79 | U | | 0.78 | U | | 0.79 | U | |
| DIBENZOFURAN | | 0.77 | U | | 0.79 | U | | 0.78 | U | | 0.79 | U | |
| DIETHYL PHTHALATE | | 19 | U | | 19 | U | | 19 | U | | 19 | U | |
| DIMETHYL PHTHALATE | | 20 | U | | 20 | U | | 20 | U | | 20 | U | |
| DI-N-BUTYL PHTHALATE | | 17 | U | | 18 | U | | 18 | U | | 18 | U | |
| DI-N-OCTYL PHTHALATE | | 9.2 | U | | 9.5 | U | | 9.4 | U | | 9.4 | U | |
| FLUORANTHENE | | 0.64 | U | | 0.66 | U | | 0.65 | U | | 0.66 | U | |
| FLUORENE | | 0.62 | U | | 0.64 | U | | 0.63 | U | | 0.63 | U | |
| HEXACHLOROBENZENE | | 2.4 | U | | 2.5 | U | | 2.5 | U | | 2.5 | U | |
| HEXACHLOROBUTADIENE | | 6.5 | U | | 6.7 | U | | 6.6 | U | | 6.7 | U | |
| HEXACHLOROCYCLOPENTADIENE | | 9.4 | UJ | C | 9.7 | U | | 9.6 | UJ | C | 9.7 | U | |
| HEXACHLOROETHANE | | 10 | U | | 11 | U | | 11 | U | | 11 | U | |
| INDENO(1,2,3-CD)PYRENE | | 0.41 | U | | 0.42 | U | | 0.42 | U | | 0.42 | U | |
| ISOPHORONE | | 15 | U | | 16 | U | | 15 | U | | 16 | U | |
| NAPHTHALENE | | 0.95 | U | | 0.99 | U | | 0.97 | U | | 0.98 | U | |
| NITROBENZENE | | 2.6 | U | | 2.6 | U | | 2.6 | U | | 2.6 | U | |
| N-NITROSO-DI-N-PROPYLAMINE | | 7.3 | U | | 7.6 | UJ | C | 7.5 | U | | 7.5 | UJ | C |
| N-NITROSODIPHENYLAMINE | | 24 | U | | 25 | U | | 25 | U | | 25 | U | |
| PENTACHLOROPHENOL | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| PHENANTHRENE | | 0.85 | U | | 0.88 | U | | 0.87 | U | | 0.87 | U | |
| PHENOL | | 8.5 | U | | 8.8 | U | | 8.7 | U | | 8.7 | U | |
| PYRENE | | 0.51 | U | | 0.53 | U | | 0.52 | U | | 0.53 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-13-14 | | | F-VB-UST-14-14 | | | F-VB-UST-2-14 | | | F-VB-UST-3-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-13 | | | 240-57597-14 | | | 240-57597-2 | | | 240-57597-3 | | |
| **FRACTION: OS** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 85.3 | | | 85.5 | | | 85.0 | | | 82.2 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| BENZO(K)FLUORANTHENE | | 0.79 | U | | 0.8 | U | | 0.8 | U | | 0.83 | U | |
| BIS(2-CHLOROETHOXY)METHANE | | 25 | U | | 26 | U | | 26 | U | | 27 | U | |
| BIS(2-CHLOROETHYL)ETHER | | 2.3 | U | | 2.4 | U | | 2.3 | U | | 2.4 | U | |
| BIS(2-ETHYLHEXYL)PHTHALATE | | 60 | J | P | 22 | U | | 22 | U | | 23 | U | |
| BUTYL BENZYL PHTHALATE | | 12 | U | | 12 | U | | 12 | U | | 12 | U | |
| CAPROLACTAM | | 43 | U | | 44 | U | | 43 | U | | 45 | U | |
| CARBAZOLE | | 31 | U | | 32 | U | | 32 | U | | 33 | U | |
| CHRYSENE | | 1.3 | U | | 1.3 | U | | 1.3 | U | | 1.3 | U | |
| DIBENZO(A,H)ANTHRACENE | | 0.76 | U | | 0.78 | U | | 0.77 | U | | 0.8 | U | |
| DIBENZOFURAN | | 0.76 | U | | 0.78 | U | | 0.77 | U | | 0.8 | U | |
| DIETHYL PHTHALATE | | 19 | U | | 19 | U | | 19 | U | | 19 | U | |
| DIMETHYL PHTHALATE | | 20 | U | | 20 | U | | 20 | U | | 21 | U | |
| DI-N-BUTYL PHTHALATE | | 18 | J | P | 18 | U | | 18 | U | | 18 | U | |
| DI-N-OCTYL PHTHALATE | | 9.1 | U | | 9.4 | U | | 9.2 | U | | 9.6 | U | |
| FLUORANTHENE | | 0.64 | U | | 0.65 | U | | 0.64 | U | | 0.67 | U | |
| FLUORENE | | 0.61 | U | | 0.63 | U | | 0.62 | U | | 0.64 | U | |
| HEXACHLOROBENZENE | | 2.4 | U | | 2.5 | U | | 2.5 | U | | 2.6 | U | |
| HEXACHLOROBUTADIENE | | 6.5 | U | | 6.6 | U | | 6.6 | U | | 6.8 | U | |
| HEXACHLOROCYCLOPENTADIENE | | 9.4 | U | | 9.6 | U | | 9.5 | UJ | C | 9.8 | UJ | C |
| HEXACHLOROETHANE | | 10 | U | | 11 | U | | 11 | U | | 11 | U | |
| INDENO(1,2,3-CD)PYRENE | | 0.41 | U | | 0.41 | U | | 0.41 | U | | 0.43 | U | |
| ISOPHORONE | | 15 | U | | 15 | U | | 15 | U | | 16 | U | |
| NAPHTHALENE | | 0.95 | U | | 0.97 | U | | 0.96 | U | | 1 | U | |
| NITROBENZENE | | 2.5 | U | | 2.6 | U | | 2.6 | U | | 2.7 | U | |
| N-NITROSO-DI-N-PROPYLAMINE | | 7.3 | UJ | C | 7.5 | UJ | C | 7.4 | U | | 7.7 | U | |
| N-NITROSODIPHENYLAMINE | | 24 | U | | 25 | U | | 25 | U | | 26 | U | |
| PENTACHLOROPHENOL | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| PHENANTHRENE | | 0.85 | U | | 0.86 | U | | 0.85 | U | | 0.89 | U | |
| PHENOL | | 8.5 | U | | 8.6 | U | | 8.5 | U | | 8.9 | U | |
| PYRENE | | 0.51 | U | | 0.52 | U | | 0.51 | U | | 0.53 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-4-14 | | | F-VB-UST-5-14 | | | F-VB-UST-6-14 | | | F-VB-UST-7-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-4 | | | 240-57597-5 | | | 240-57597-6 | | | 240-57597-7 | | |
| **FRACTION: OS** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 84.4 | | | 84.8 | | | 83.5 | | | 83.8 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| BENZO(K)FLUORANTHENE | | 0.81 | U | | 0.79 | U | | 0.81 | U | | 0.82 | U | |
| BIS(2-CHLOROETHOXY)METHANE | | 26 | U | | 26 | U | | 26 | U | | 27 | U | |
| BIS(2-CHLOROETHYL)ETHER | | 2.4 | U | | 2.3 | U | | 2.4 | U | | 2.4 | U | |
| BIS(2-ETHYLHEXYL)PHTHALATE | | 33 | J | P | 38 | J | P | 32 | J | P | 32 | J | P |
| BUTYL BENZYL PHTHALATE | | 12 | U | | 12 | U | | 12 | U | | 12 | U | |
| CAPROLACTAM | | 44 | U | | 43 | U | | 44 | U | | 45 | U | |
| CARBAZOLE | | 32 | U | | 32 | U | | 32 | U | | 33 | U | |
| CHRYSENE | | 1.3 | U | | 1.3 | U | | 1.3 | U | | 1.3 | U | |
| DIBENZO(A,H)ANTHRACENE | | 0.79 | U | | 0.77 | U | | 0.79 | U | | 0.8 | U | |
| DIBENZOFURAN | | 0.79 | U | | 0.77 | U | | 0.79 | U | | 0.8 | U | |
| DIETHYL PHTHALATE | | 19 | U | | 19 | U | | 19 | U | | 19 | U | |
| DIMETHYL PHTHALATE | | 20 | U | | 20 | U | | 20 | U | | 21 | U | |
| DI-N-BUTYL PHTHALATE | | 18 | U | | 18 | U | | 18 | U | | 18 | U | |
| DI-N-OCTYL PHTHALATE | | 9.4 | U | | 9.2 | U | | 9.4 | U | | 9.6 | U | |
| FLUORANTHENE | | 0.66 | U | | 0.64 | U | | 0.66 | U | | 0.67 | U | |
| FLUORENE | | 0.63 | U | | 0.62 | U | | 0.63 | U | | 0.64 | U | |
| HEXACHLOROBENZENE | | 2.5 | U | | 2.5 | U | | 2.5 | U | | 2.5 | U | |
| HEXACHLOROBUTADIENE | | 6.7 | U | | 6.5 | U | | 6.7 | U | | 6.8 | U | |
| HEXACHLOROCYCLOPENTADIENE | | 9.7 | UJ | C | 9.5 | UJ | C | 9.7 | UJ | C | 9.8 | UJ | C |
| HEXACHLOROETHANE | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| INDENO(1,2,3-CD)PYRENE | | 0.42 | U | | 0.41 | U | | 0.42 | U | | 0.42 | U | |
| ISOPHORONE | | 15 | U | | 15 | U | | 15 | U | | 16 | U | |
| NAPHTHALENE | | 0.98 | U | | 0.96 | U | | 0.98 | U | | 0.99 | U | |
| NITROBENZENE | | 2.6 | U | | 2.6 | U | | 2.6 | U | | 2.7 | U | |
| N-NITROSO-DI-N-PROPYLAMINE | | 7.5 | U | | 7.4 | U | | 7.5 | U | | 7.6 | U | |
| N-NITROSODIPHENYLAMINE | | 25 | U | | 25 | U | | 25 | U | | 25 | U | |
| PENTACHLOROPHENOL | | 11 | U | | 11 | U | | 11 | U | | 11 | U | |
| PHENANTHRENE | | 0.87 | U | | 0.85 | U | | 0.87 | U | | 0.89 | U | |
| PHENOL | | 8.7 | U | | 8.5 | U | | 8.7 | U | | 8.9 | U | |
| PYRENE | | 0.52 | U | | 0.51 | U | | 0.52 | U | | 0.53 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-8-14 | | | F-VB-UST-9-14 | | | F-VB-UST-DUP | | |
|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-8 | | | 240-57597-9 | | | 240-57597-15 | | |
| **FRACTION: OS** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 82.0 | | | 80.9 | | | 85.5 | | |
| | DUP_OF | | | | | | | F-VB-UST-2-14 | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| BENZO(K)FLUORANTHENE | | 0.82 | U | | 0.85 | U | | 0.79 | U | |
| BIS(2-CHLOROETHOXY)METHANE | | 27 | U | | 27 | U | | 26 | U | |
| BIS(2-CHLOROETHYL)ETHER | | 2.4 | U | | 2.5 | U | | 2.3 | U | |
| BIS(2-ETHYLHEXYL)PHTHALATE | | 23 | U | | 40 | J | P | 22 | J | P |
| BUTYL BENZYL PHTHALATE | | 12 | U | | 12 | U | | 12 | U | |
| CAPROLACTAM | | 45 | U | | 46 | U | | 43 | U | |
| CARBAZOLE | | 33 | U | | 34 | U | | 31 | U | |
| CHRYSENE | | 1.3 | U | | 1.4 | U | | 1.3 | U | |
| DIBENZO(A,H)ANTHRACENE | | 0.8 | U | | 0.82 | U | | 0.77 | U | |
| DIBENZOFURAN | | 0.8 | U | | 0.82 | U | | 0.77 | U | |
| DIETHYL PHTHALATE | | 19 | U | | 20 | U | | 19 | U | |
| DIMETHYL PHTHALATE | | 21 | U | | 21 | U | | 20 | U | |
| DI-N-BUTYL PHTHALATE | | 18 | U | | 19 | U | | 17 | U | |
| DI-N-OCTYL PHTHALATE | | 9.6 | U | | 9.9 | U | | 9.2 | U | |
| FLUORANTHENE | | 0.67 | U | | 0.69 | U | | 0.64 | U | |
| FLUORENE | | 0.64 | U | | 0.66 | U | | 0.62 | U | |
| HEXACHLOROBENZENE | | 2.5 | U | | 2.6 | U | | 2.4 | U | |
| HEXACHLOROBUTADIENE | | 6.8 | U | | 7 | U | | 6.5 | U | |
| HEXACHLOROCYCLOPENTADIENE | | 9.8 | UJ | C | 10 | UJ | C | 9.4 | U | |
| HEXACHLOROETHANE | | 11 | U | | 11 | U | | 10 | U | |
| INDENO(1,2,3-CD)PYRENE | | 0.42 | U | | 0.44 | U | | 0.41 | U | |
| ISOPHORONE | | 16 | U | | 16 | U | | 15 | U | |
| NAPHTHALENE | | 0.99 | U | | 1 | U | | 0.96 | U | |
| NITROBENZENE | | 2.7 | U | | 2.7 | U | | 2.6 | U | |
| N-NITROSO-DI-N-PROPYLAMINE | | 7.6 | U | | 7.9 | U | | 7.3 | UJ | C |
| N-NITROSODIPHENYLAMINE | | 25 | U | | 26 | U | | 24 | U | |
| PENTACHLOROPHENOL | | 11 | U | | 11 | U | | 11 | U | |
| PHENANTHRENE | | 0.88 | U | | 0.91 | U | | 0.85 | U | |
| PHENOL | | 8.8 | U | | 9.1 | U | | 8.5 | U | |
| PYRENE | | 0.53 | U | | 0.55 | U | | 0.51 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-10-14 | | | F-VB-UST-11-14 | | | F-VB-UST-1-14 | | | F-VB-UST-12-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-10 | | | 240-57597-11 | | | 240-57597-1 | | | 240-57597-12 | | |
| **FRACTION: PCB** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 86.6 | | | 83.5 | | | 84.3 | | | 84.1 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| AROCLOR-1016 | | 14 | U | | 14 | U | | 14 | U | | 14 | U | |
| AROCLOR-1221 | | 19 | U | | 19 | U | | 19 | U | | 19 | U | |
| AROCLOR-1232 | | 23 | U | | 24 | U | | 24 | U | | 24 | U | |
| AROCLOR-1242 | | 13 | U | | 13 | U | | 13 | U | | 13 | U | |
| AROCLOR-1248 | | 9.3 | U | | 9.5 | U | | 9.6 | U | | 9.4 | U | |
| AROCLOR-1254 | | 16 | U | | 17 | U | | 17 | U | | 16 | U | |
| AROCLOR-1260 | | 10 | U | | 11 | U | | 11 | U | | 11 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-13-14 | | | F-VB-UST-14-14 | | | F-VB-UST-2-14 | | | F-VB-UST-3-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-13 | | | 240-57597-14 | | | 240-57597-2 | | | 240-57597-3 | | |
| **FRACTION: PCB** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 85.3 | | | 85.5 | | | 85.0 | | | 82.2 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| AROCLOR-1016 | | 14 | U | | 14 | U | | 14 | U | | 15 | U | |
| AROCLOR-1221 | | 19 | U | | 19 | U | | 19 | U | | 19 | U | |
| AROCLOR-1232 | | 23 | U | | 23 | U | | 23 | U | | 24 | U | |
| AROCLOR-1242 | | 13 | U | | 13 | U | | 13 | U | | 13 | U | |
| AROCLOR-1248 | | 9.3 | U | | 9.3 | U | | 9.4 | U | | 9.7 | U | |
| AROCLOR-1254 | | 16 | U | | 16 | U | | 16 | U | | 17 | U | |
| AROCLOR-1260 | | 10 | U | | 10 | U | | 11 | U | | 11 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-4-14 | | | F-VB-UST-5-14 | | | F-VB-UST-6-14 | | | F-VB-UST-7-14 | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-4 | | | 240-57597-5 | | | 240-57597-6 | | | 240-57597-7 | | |
| **FRACTION: PCB** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 84.4 | | | 84.8 | | | 83.5 | | | 83.8 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| AROCLOR-1016 | | 14 | U | | 14 | U | | 14 | U | | 14 | U | |
| AROCLOR-1221 | | 19 | U | | 19 | U | | 19 | U | | 19 | U | |
| AROCLOR-1232 | | 24 | U | | 23 | U | | 24 | U | | 24 | U | |
| AROCLOR-1242 | | 13 | U | | 13 | U | | 13 | U | | 13 | U | |
| AROCLOR-1248 | | 9.6 | U | | 9.3 | U | | 9.6 | U | | 9.5 | U | |
| AROCLOR-1254 | | 17 | U | | 16 | U | | 17 | U | | 17 | U | |
| AROCLOR-1260 | | 11 | U | | 10 | U | | 11 | U | | 11 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-8-14 | | | F-VB-UST-9-14 | | | F-VB-UST-DUP | | |
|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-8 | | | 240-57597-9 | | | 240-57597-15 | | |
| **FRACTION: PCB** | SAMP_DATE | 11/5/2015 | | | 11/5/2015 | | | 11/5/2015 | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | NM | | | NM | | |
| | UNITS | UG/KG | | | UG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 82.0 | | | 80.9 | | | 85.5 | | |
| | DUP_OF | | | | | | | F-VB-UST-2-14 | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| AROCLOR-1016 | | 15 | U | | 15 | U | | 14 | U | |
| AROCLOR-1221 | | 20 | U | | 20 | U | | 18 | U | |
| AROCLOR-1232 | | 25 | U | | 25 | U | | 23 | U | |
| AROCLOR-1242 | | 14 | U | | 14 | U | | 13 | U | |
| AROCLOR-1248 | | 9.8 | U | | 9.9 | U | | 9.2 | U | |
| AROCLOR-1254 | | 17 | U | | 17 | U | | 16 | U | |
| AROCLOR-1260 | | 11 | U | | 11 | U | | 10 | U | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-10-14 | | | | | | F-VB-UST-11-14 | | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-10 | | | | | | 240-57597-11 | | | | | |
| **FRACTION: PET** | SAMP_DATE | 11/5/2015 | | | | | | 11/5/2015 | | | | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | | | | NM | | | | | |
| | UNITS | MG/KG | | | UG/KG | | | MG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 86.6 | | | 86.6 | | | 83.5 | | | 83.5 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| TPH (C06-C10) | | | | | 53 | U | | | | | 55 | U | |
| TPH (C10-C32) | | 13 | U | A | | | | 9 | U | A | | | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-1-14 | | | | | | F-VB-UST-12-14 | | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-1 | | | | | | 240-57597-12 | | | | | |
| **FRACTION: PET** | SAMP_DATE | 11/5/2015 | | | | | | 11/5/2015 | | | | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | | | | NM | | | | | |
| | UNITS | MG/KG | | | UG/KG | | | MG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 84.3 | | | 84.3 | | | 84.1 | | | 84.1 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| TPH (C06-C10) | | | | | 55 | U | | | | | 54 | U | |
| TPH (C10-C32) | | 16 | U | A | | | | 30 | U | A | | | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-13-14 | | | | | | F-VB-UST-14-14 | | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-13 | | | | | | 240-57597-14 | | | | | |
| **FRACTION: PET** | SAMP_DATE | 11/5/2015 | | | | | | 11/5/2015 | | | | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | | | | NM | | | | | |
| | UNITS | MG/KG | | | UG/KG | | | MG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 85.3 | | | 85.3 | | | 85.5 | | | 85.5 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| TPH (C06-C10) | | | | | 54 | U | | | | | 54 | U | |
| TPH (C10-C32) | | 13 | U | A | | | | 11 | U | A | | | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-2-14 | | | | | | F-VB-UST-3-14 | | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-2 | | | | | | 240-57597-3 | | | | | |
| **FRACTION: PET** | SAMP_DATE | 11/5/2015 | | | | | | 11/5/2015 | | | | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | | | | NM | | | | | |
| | UNITS | MG/KG | | | UG/KG | | | MG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 85.0 | | | 85.0 | | | 82.2 | | | 82.2 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| TPH (C06-C10) | | | | | 54 | U | | | | | 56 | U | |
| TPH (C10-C32) | | 10 | U | A | | | | 7.7 | U | | | | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-4-14 | | | | | | F-VB-UST-5-14 | | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-4 | | | | | | 240-57597-5 | | | | | |
| **FRACTION: PET** | SAMP_DATE | 11/5/2015 | | | | | | 11/5/2015 | | | | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | | | | NM | | | | | |
| | UNITS | MG/KG | | | UG/KG | | | MG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 84.4 | | | 84.4 | | | 84.8 | | | 84.8 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| TPH (C06-C10) | | | | | 55 | U | | | | | 77 | J | P |
| TPH (C10-C32) | | 7.6 | U | A | | | | 7.5 | U | | | | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-6-14 | | | | | | F-VB-UST-7-14 | | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-6 | | | | | | 240-57597-7 | | | | | |
| **FRACTION: PET** | SAMP_DATE | 11/5/2015 | | | | | | 11/5/2015 | | | | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | | | | NM | | | | | |
| | UNITS | MG/KG | | | UG/KG | | | MG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 83.5 | | | 83.5 | | | 83.8 | | | 83.8 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| TPH (C06-C10) | | | | | 55 | U | | | | | 55 | U | |
| TPH (C10-C32) | | 7.7 | U | | | | | 16 | U | A | | | |

| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-8-14 | | | | | | F-VB-UST-9-14 | | | | | |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| **SDG: 240-57597-1** | LAB_ID | 240-57597-8 | | | | | | 240-57597-9 | | | | | |
| **FRACTION: PET** | SAMP_DATE | 11/5/2015 | | | | | | 11/5/2015 | | | | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | | | | NM | | | | | |
| | UNITS | MG/KG | | | UG/KG | | | MG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 82.0 | | | 82.0 | | | 80.9 | | | 80.9 | | |
| | DUP_OF | | | | | | | | | | | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| TPH (C06-C10) | | | | | 56 | U | | | | | 57 | U | |
| TPH (C10-C32) | | 7.7 | U | | | | | 7.8 | U | | | | |

| | | | | | | | |
|---|---|---|---|---|---|---|---|
| **PROJ_NO: 07336** | NSAMPLE | F-VB-UST-DUP | | | | | |
| **SDG: 240-57597-1** | LAB_ID | 240-57597-15 | | | | | |
| **FRACTION: PET** | SAMP_DATE | 11/5/2015 | | | | | |
| **MEDIA: SOIL** | QC_TYPE | NM | | | | | |
| | UNITS | MG/KG | | | UG/KG | | |
| | PCT_SOLIDS | 85.5 | | | 85.5 | | |
| | DUP_OF | F-VB-UST-2-14 | | | F-VB-UST-2-14 | | |
| PARAMETER | | RESULT | VQL | QLCD | RESULT | VQL | QLCD |
| TPH (C06-C10) | | | | | 54 | U | |
| TPH (C10-C32) | | 8.2 | U | A | | | |

**Appendix B**

Results as Reported by the Laboratory

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 146336.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 5.89(g) | Date Analyzed: 11/10/2015 14:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.81 | U | 5.0 | 0.81 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.79 | U | 5.0 | 0.79 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.91 | U | 5.0 | 0.91 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 146336.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 5.89(g) | Date Analyzed: 11/10/2015 14:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.80 | U | 5.0 | 0.80 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.72 | U | 5.0 | 0.72 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.8 | U | 200 | 5.8 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.75 | U | 5.0 | 0.75 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.35 | U | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.92 | U | 5.0 | 0.92 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.81 | U | 5.0 | 0.81 |
| 108-88-3 | Toluene | 0.27 | U | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 146336.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 5.89(g) | Date Analyzed: 11/10/2015 14:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 103 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 91 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 103 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 146337.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 5.36(g) | Date Analyzed: 11/10/2015 15:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.71 | U | 5.5 | 0.71 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.37 | U | 5.5 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 0.88 | U | 5.5 | 0.88 |
| 563-58-6 | 1,1-Dichloropropene | 0.41 | U | 5.5 | 0.41 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.86 | U | 5.5 | 0.86 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.50 | U | 5.5 | 0.50 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.42 | U | 5.5 | 0.42 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.64 | U | 5.5 | 0.64 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.37 | U | 5.5 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 0.50 | U | 5.5 | 0.50 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.76 | U | 5.5 | 0.76 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.48 | U | 5.5 | 0.48 |
| 78-93-3 | 2-Butanone (MEK) | 4.0 | J B | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.70 | U | 55 | 0.70 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.75 | U | 22 | 0.75 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.5 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.45 | U | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 22 | 1.3 |
| 67-64-1 | Acetone | 6.3 | U | 22 | 6.3 |
| 71-43-2 | Benzene | 0.99 | U | 5.5 | 0.99 |
| 108-86-1 | Bromobenzene | 0.36 | U | 5.5 | 0.36 |
| 75-25-2 | Bromoform | 0.26 | U | 5.5 | 0.26 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.65 | U | 5.5 | 0.65 |
| 56-23-5 | Carbon tetrachloride | 0.71 | U | 5.5 | 0.71 |
| 108-90-7 | Chlorobenzene | 0.58 | U | 5.5 | 0.58 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 146337.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 5.36(g) | Date Analyzed: 11/10/2015 15:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.41 | U | 5.5 | 0.41 |
| 75-00-3 | Chloroethane | 0.48 | U | 5.5 | 0.48 |
| 67-66-3 | Chloroform | 0.39 | U | 5.5 | 0.39 |
| 74-87-3 | Chloromethane | 0.87 | U | 5.5 | 0.87 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.78 | U | 5.5 | 0.78 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.36 | U | 5.5 | 0.36 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.3 | U | 220 | 6.3 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.65 | U | 11 | 0.65 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.47 | U | 5.5 | 0.47 |
| 75-09-2 | Methylene Chloride | 0.81 | U | 5.5 | 0.81 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.38 | U | 5.5 | 0.38 |
| 104-51-8 | n-Butylbenzene | 0.65 | U | 5.5 | 0.65 |
| 103-65-1 | N-Propylbenzene | 0.66 | U | 5.5 | 0.66 |
| 95-47-6 | o-Xylene | 0.63 | U | 5.5 | 0.63 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.5 | 1.0 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.47 | U | 5.5 | 0.47 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.88 | U | 5.5 | 0.88 |
| 108-88-3 | Toluene | 0.30 | U | 5.5 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.42 | U | 5.5 | 0.42 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.72 | U | 11 | 0.72 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.59 | U | 11 | 0.59 |
| 98-06-6 | tert-Butylbenzene | 0.74 | U | 5.5 | 0.74 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 146337.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 5.36(g) | Date Analyzed: 11/10/2015 15:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.58 | U | 5.5 | 0.58 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 96 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 95 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 85 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 99 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 146338.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 5.23(g) | Date Analyzed: 11/10/2015 15:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.76 | U | 5.8 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 5.8 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.91 | U | 5.8 | 0.91 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 5.8 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 5.8 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.48 | U | 5.8 | 0.48 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 6.6 | U | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.69 | U | 5.8 | 0.69 |
| 56-23-5 | Carbon tetrachloride | 0.76 | U | 5.8 | 0.76 |
| 108-90-7 | Chlorobenzene | 0.62 | U | 5.8 | 0.62 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 146338.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 5.23(g) | Date Analyzed: 11/10/2015 15:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 5.8 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.83 | U | 5.8 | 0.83 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 5.8 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.69 | U | 12 | 0.69 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.41 | U | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 0.69 | U | 5.8 | 0.69 |
| 103-65-1 | N-Propylbenzene | 0.70 | U | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.48 | U | 5.8 | 0.48 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.77 | U | 12 | 0.77 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.63 | U | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 0.78 | U | 5.8 | 0.78 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 146338.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 5.23(g) | Date Analyzed: 11/10/2015 15:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.62 | U | 5.8 | 0.62 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 93 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 90 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 86 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 90 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 146339.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 5.43(g) | Date Analyzed: 11/10/2015 15:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.71 | U | 5.5 | 0.71 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.37 | U | 5.5 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 0.87 | U | 5.5 | 0.87 |
| 563-58-6 | 1,1-Dichloropropene | 0.40 | U | 5.5 | 0.40 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.85 | U | 5.5 | 0.85 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.50 | U | 5.5 | 0.50 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.41 | U | 5.5 | 0.41 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.63 | U | 5.5 | 0.63 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.37 | U | 5.5 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 0.50 | U | 5.5 | 0.50 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 5.5 | 0.75 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.48 | U | 5.5 | 0.48 |
| 78-93-3 | 2-Butanone (MEK) | 4.0 | J B | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.70 | U | 55 | 0.70 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.74 | U | 22 | 0.74 |
| 106-43-4 | 4-Chlorotoluene | 0.28 | U | 5.5 | 0.28 |
| 99-87-6 | p-Isopropyltoluene | 0.45 | U | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 22 | 1.3 |
| 67-64-1 | Acetone | 6.2 | U | 22 | 6.2 |
| 71-43-2 | Benzene | 0.98 | U | 5.5 | 0.98 |
| 108-86-1 | Bromobenzene | 0.36 | U | 5.5 | 0.36 |
| 75-25-2 | Bromoform | 0.26 | U | 5.5 | 0.26 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.64 | U | 5.5 | 0.64 |
| 56-23-5 | Carbon tetrachloride | 0.71 | U | 5.5 | 0.71 |
| 108-90-7 | Chlorobenzene | 0.58 | U | 5.5 | 0.58 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 146339.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 5.43(g) | Date Analyzed: 11/10/2015 15:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.40 | U | 5.5 | 0.40 |
| 75-00-3 | Chloroethane | 0.48 | U | 5.5 | 0.48 |
| 67-66-3 | Chloroform | 0.39 | U | 5.5 | 0.39 |
| 74-87-3 | Chloromethane | 0.86 | U | 5.5 | 0.86 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.77 | U | 5.5 | 0.77 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.36 | U | 5.5 | 0.36 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.2 | U | 220 | 6.2 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.64 | U | 11 | 0.64 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.47 | U | 5.5 | 0.47 |
| 75-09-2 | Methylene Chloride | 0.81 | U | 5.5 | 0.81 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.38 | U | 5.5 | 0.38 |
| 104-51-8 | n-Butylbenzene | 0.64 | U | 5.5 | 0.64 |
| 103-65-1 | N-Propylbenzene | 0.65 | U | 5.5 | 0.65 |
| 95-47-6 | o-Xylene | 0.62 | U | 5.5 | 0.62 |
| 135-98-8 | sec-Butylbenzene | 0.99 | U | 5.5 | 0.99 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.47 | U | 5.5 | 0.47 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.87 | U | 5.5 | 0.87 |
| 108-88-3 | Toluene | 0.29 | U | 5.5 | 0.29 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.41 | U | 5.5 | 0.41 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.72 | U | 11 | 0.72 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.59 | U | 11 | 0.59 |
| 98-06-6 | tert-Butylbenzene | 0.73 | U | 5.5 | 0.73 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 146339.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 5.43(g) | Date Analyzed: 11/10/2015 15:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.58 | U | 5.5 | 0.58 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 93 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 94 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 83 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 146340.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.09(g) | Date Analyzed: 11/10/2015 16:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U F2 | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.3 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U F2 | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U F1 | 23 | 1.4 |
| 67-64-1 | Acetone | 13 | J | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 146340.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.09(g) | Date Analyzed: 11/10/2015 16:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U F2 | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.41 | J | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 0.68 | U F2 | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.70 | U | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U F2 | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U F1 | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.63 | U | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 0.78 | U | 5.8 | 0.78 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 146340.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.09(g) | Date Analyzed: 11/10/2015 16:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 100 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 104 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 146343.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.18(g) | Date Analyzed: 11/10/2015 17:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 3.8 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 6.6 | U | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 146343.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.18(g) | Date Analyzed: 11/10/2015 17:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.45 | J | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 146343.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.18(g) | Date Analyzed: 11/10/2015 17:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 88 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 89 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 81 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 146344.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 5.31(g) | Date Analyzed: 11/10/2015 17:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.56 | U | 5.6 | 0.56 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.73 | U | 5.6 | 0.73 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.35 | U | 5.6 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.47 | U | 5.6 | 0.47 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.6 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.90 | U | 5.6 | 0.90 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 5.6 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.56 | U | 5.6 | 0.56 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.88 | U | 5.6 | 0.88 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.52 | U | 5.6 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.43 | U | 5.6 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.65 | U | 5.6 | 0.65 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.6 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.52 | U | 5.6 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.6 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.78 | U | 5.6 | 0.78 |
| 142-28-9 | 1,3-Dichloropropane | 0.45 | U | 5.6 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.6 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.6 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 4.1 | J B | 22 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.72 | U | 56 | 0.72 |
| 95-49-8 | 2-Chlorotoluene | 0.34 | U | 5.6 | 0.34 |
| 591-78-6 | 2-Hexanone | 0.76 | U | 22 | 0.76 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.6 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.46 | U | 5.6 | 0.46 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 6.4 | U | 22 | 6.4 |
| 71-43-2 | Benzene | 1.0 | U | 5.6 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.6 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.6 | 0.27 |
| 74-83-9 | Bromomethane | 0.47 | U | 5.6 | 0.47 |
| 75-15-0 | Carbon disulfide | 0.66 | U | 5.6 | 0.66 |
| 56-23-5 | Carbon tetrachloride | 0.73 | U | 5.6 | 0.73 |
| 108-90-7 | Chlorobenzene | 0.60 | U | 5.6 | 0.60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 146344.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 5.31(g) | Date Analyzed: 11/10/2015 17:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.42 | U | 5.6 | 0.42 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.6 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.6 | 0.40 |
| 74-87-3 | Chloromethane | 0.89 | U | 5.6 | 0.89 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.6 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.80 | U | 5.6 | 0.80 |
| 74-95-3 | Dibromomethane | 0.47 | U | 5.6 | 0.47 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.6 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.6 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.6 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.4 | U | 220 | 6.4 |
| 100-41-4 | Ethylbenzene | 0.36 | U | 5.6 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 0.45 | U | 5.6 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.66 | U | 11 | 0.66 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.6 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.6 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.83 | U | 5.6 | 0.83 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.47 | U | 11 | 0.47 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.6 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.66 | U | 5.6 | 0.66 |
| 103-65-1 | N-Propylbenzene | 0.67 | U | 5.6 | 0.67 |
| 95-47-6 | o-Xylene | 0.64 | U | 5.6 | 0.64 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.6 | 1.0 |
| 100-42-5 | Styrene | 0.45 | U | 5.6 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.6 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.46 | U | 5.6 | 0.46 |
| 127-18-4 | Tetrachloroethene | 0.90 | U | 5.6 | 0.90 |
| 108-88-3 | Toluene | 0.30 | U | 5.6 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.36 | U | 5.6 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.6 | 0.33 |
| 79-01-6 | Trichloroethene | 0.43 | U | 5.6 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.6 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.74 | U | 11 | 0.74 |
| 75-01-4 | Vinyl chloride | 0.34 | U | 5.6 | 0.34 |
| 1330-20-7 | Xylenes, Total | 0.61 | U | 11 | 0.61 |
| 98-06-6 | tert-Butylbenzene | 0.75 | U | 5.6 | 0.75 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 146344.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 5.31(g) | Date Analyzed: 11/10/2015 17:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.60 | U | 5.6 | 0.60 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 98 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 103 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 146345.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 5.29(g) | Date Analyzed: 11/10/2015 17:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.48 | U | 5.8 | 0.48 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.92 | U | 5.8 | 0.92 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.3 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.78 | U | 23 | 0.78 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 6.6 | U | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.48 | U | 5.8 | 0.48 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 146345.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 5.29(g) | Date Analyzed: 11/10/2015 17:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.41 | U | 5.8 | 0.41 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.48 | U | 5.8 | 0.48 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.8 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.85 | U | 5.8 | 0.85 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.48 | U | 12 | 0.48 |
| 91-20-3 | Naphthalene | 0.40 | U | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.8 | 1.0 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.92 | U | 5.8 | 0.92 |
| 108-88-3 | Toluene | 0.33 | J | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.8 | 0.33 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.8 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 146345.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 5.29(g) | Date Analyzed: 11/10/2015 17:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 97 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 93 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 85 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 146346.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.35(g) | Date Analyzed: 11/10/2015 18:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.92 | U | 5.8 | 0.92 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.4 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 12 | J | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 146346.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.35(g) | Date Analyzed: 11/10/2015 18:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.85 | U | 5.8 | 0.85 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.40 | U | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.92 | U | 5.8 | 0.92 |
| 108-88-3 | Toluene | 0.36 | J | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 146346.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.35(g) | Date Analyzed: 11/10/2015 18:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 88 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 82 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 146347.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.14(g) | Date Analyzed: 11/10/2015 18:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.56 | U | 5.6 | 0.56 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.73 | U | 5.6 | 0.73 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.35 | U | 5.6 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.47 | U | 5.6 | 0.47 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.6 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.90 | U | 5.6 | 0.90 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 5.6 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.56 | U | 5.6 | 0.56 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.88 | U | 5.6 | 0.88 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.52 | U | 5.6 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.43 | U | 5.6 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.65 | U | 5.6 | 0.65 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.6 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.52 | U | 5.6 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.6 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.78 | U | 5.6 | 0.78 |
| 142-28-9 | 1,3-Dichloropropane | 0.45 | U | 5.6 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.6 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.6 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 3.5 | J B | 22 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.72 | U | 56 | 0.72 |
| 95-49-8 | 2-Chlorotoluene | 0.34 | U | 5.6 | 0.34 |
| 591-78-6 | 2-Hexanone | 0.76 | U | 22 | 0.76 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.6 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.46 | U | 5.6 | 0.46 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 6.4 | U | 22 | 6.4 |
| 71-43-2 | Benzene | 1.0 | U | 5.6 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.6 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.6 | 0.27 |
| 74-83-9 | Bromomethane | 0.47 | U | 5.6 | 0.47 |
| 75-15-0 | Carbon disulfide | 0.66 | U | 5.6 | 0.66 |
| 56-23-5 | Carbon tetrachloride | 0.73 | U | 5.6 | 0.73 |
| 108-90-7 | Chlorobenzene | 0.60 | U | 5.6 | 0.60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 146347.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.14(g) | Date Analyzed: 11/10/2015 18:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.42 | U | 5.6 | 0.42 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.6 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.6 | 0.40 |
| 74-87-3 | Chloromethane | 0.89 | U | 5.6 | 0.89 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.6 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.80 | U | 5.6 | 0.80 |
| 74-95-3 | Dibromomethane | 0.47 | U | 5.6 | 0.47 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.6 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.6 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.6 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.4 | U | 220 | 6.4 |
| 100-41-4 | Ethylbenzene | 0.36 | U | 5.6 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 0.45 | U | 5.6 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.66 | U | 11 | 0.66 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.6 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.6 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.83 | U | 5.6 | 0.83 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.47 | U | 11 | 0.47 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.6 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.66 | U | 5.6 | 0.66 |
| 103-65-1 | N-Propylbenzene | 0.67 | U | 5.6 | 0.67 |
| 95-47-6 | o-Xylene | 0.64 | U | 5.6 | 0.64 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.6 | 1.0 |
| 100-42-5 | Styrene | 0.45 | U | 5.6 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.6 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.46 | U | 5.6 | 0.46 |
| 127-18-4 | Tetrachloroethene | 0.90 | U | 5.6 | 0.90 |
| 108-88-3 | Toluene | 0.30 | U | 5.6 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.36 | U | 5.6 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.6 | 0.33 |
| 79-01-6 | Trichloroethene | 0.43 | U | 5.6 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.6 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.74 | U | 11 | 0.74 |
| 75-01-4 | Vinyl chloride | 0.34 | U | 5.6 | 0.34 |
| 1330-20-7 | Xylenes, Total | 0.61 | U | 11 | 0.61 |
| 98-06-6 | tert-Butylbenzene | 0.75 | U | 5.6 | 0.75 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 146347.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.14(g) | Date Analyzed: 11/10/2015 18:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.60 | U | 5.6 | 0.60 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 88 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 83 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 146348.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/10/2015 19:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.60 | U | 6.0 | 0.60 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.77 | U | 6.0 | 0.77 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.37 | U | 6.0 | 0.37 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.50 | U | 6.0 | 0.50 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 6.0 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.95 | U | 6.0 | 0.95 |
| 563-58-6 | 1,1-Dichloropropene | 0.44 | U | 6.0 | 0.44 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.60 | U | 6.0 | 0.60 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.93 | U | 6.0 | 0.93 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.55 | U | 6.0 | 0.55 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.45 | U | 6.0 | 0.45 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.69 | U | 6.0 | 0.69 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 6.0 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.55 | U | 6.0 | 0.55 |
| 78-87-5 | 1,2-Dichloropropane | 0.17 | U | 6.0 | 0.17 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.82 | U | 6.0 | 0.82 |
| 142-28-9 | 1,3-Dichloropropane | 0.48 | U | 6.0 | 0.48 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 6.0 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.52 | U | 6.0 | 0.52 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 24 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.76 | U | 60 | 0.76 |
| 95-49-8 | 2-Chlorotoluene | 0.36 | U | 6.0 | 0.36 |
| 591-78-6 | 2-Hexanone | 0.81 | U | 24 | 0.81 |
| 106-43-4 | 4-Chlorotoluene | 0.31 | U | 6.0 | 0.31 |
| 99-87-6 | p-Isopropyltoluene | 0.49 | U | 6.0 | 0.49 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.5 | U | 24 | 1.5 |
| 67-64-1 | Acetone | 6.8 | U | 24 | 6.8 |
| 71-43-2 | Benzene | 1.1 | U | 6.0 | 1.1 |
| 108-86-1 | Bromobenzene | 0.39 | U | 6.0 | 0.39 |
| 75-25-2 | Bromoform | 0.29 | U | 6.0 | 0.29 |
| 74-83-9 | Bromomethane | 0.50 | U | 6.0 | 0.50 |
| 75-15-0 | Carbon disulfide | 0.70 | U | 6.0 | 0.70 |
| 56-23-5 | Carbon tetrachloride | 0.77 | U | 6.0 | 0.77 |
| 108-90-7 | Chlorobenzene | 0.63 | U | 6.0 | 0.63 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 146348.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/10/2015 19:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.60 | U | 6.0 | 0.60 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.77 | U | 6.0 | 0.77 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.37 | U | 6.0 | 0.37 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.50 | U | 6.0 | 0.50 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 6.0 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.95 | U | 6.0 | 0.95 |
| 563-58-6 | 1,1-Dichloropropene | 0.44 | U | 6.0 | 0.44 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.60 | U | 6.0 | 0.60 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.93 | U | 6.0 | 0.93 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.55 | U | 6.0 | 0.55 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.45 | U | 6.0 | 0.45 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.69 | U | 6.0 | 0.69 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 6.0 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.55 | U | 6.0 | 0.55 |
| 78-87-5 | 1,2-Dichloropropane | 0.17 | U | 6.0 | 0.17 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.82 | U | 6.0 | 0.82 |
| 142-28-9 | 1,3-Dichloropropane | 0.48 | U | 6.0 | 0.48 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 6.0 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.52 | U | 6.0 | 0.52 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 24 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.76 | U | 60 | 0.76 |
| 95-49-8 | 2-Chlorotoluene | 0.36 | U | 6.0 | 0.36 |
| 591-78-6 | 2-Hexanone | 0.81 | U | 24 | 0.81 |
| 106-43-4 | 4-Chlorotoluene | 0.31 | U | 6.0 | 0.31 |
| 99-87-6 | p-Isopropyltoluene | 0.49 | U | 6.0 | 0.49 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.5 | U | 24 | 1.5 |
| 67-64-1 | Acetone | 6.8 | U | 24 | 6.8 |
| 71-43-2 | Benzene | 1.1 | U | 6.0 | 1.1 |
| 108-86-1 | Bromobenzene | 0.39 | U | 6.0 | 0.39 |
| 75-25-2 | Bromoform | 0.29 | U | 6.0 | 0.29 |
| 74-83-9 | Bromomethane | 0.50 | U | 6.0 | 0.50 |
| 75-15-0 | Carbon disulfide | 0.70 | U | 6.0 | 0.70 |
| 56-23-5 | Carbon tetrachloride | 0.77 | U | 6.0 | 0.77 |
| 108-90-7 | Chlorobenzene | 0.63 | U | 6.0 | 0.63 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 146348.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/10/2015 19:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.44 | U | 6.0 | 0.44 |
| 75-00-3 | Chloroethane | 0.52 | U | 6.0 | 0.52 |
| 67-66-3 | Chloroform | 0.43 | U | 6.0 | 0.43 |
| 74-87-3 | Chloromethane | 0.94 | U | 6.0 | 0.94 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 6.0 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.85 | U | 6.0 | 0.85 |
| 74-95-3 | Dibromomethane | 0.50 | U | 6.0 | 0.50 |
| 74-97-5 | Bromochloromethane | 0.39 | U | 6.0 | 0.39 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 6.0 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.35 | U | 6.0 | 0.35 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.8 | U | 240 | 6.8 |
| 100-41-4 | Ethylbenzene | 0.38 | U | 6.0 | 0.38 |
| 87-68-3 | Hexachlorobutadiene | 0.48 | U | 6.0 | 0.48 |
| 108-20-3 | Diisopropyl ether | 0.70 | U | 12 | 0.70 |
| 98-82-8 | Isopropylbenzene | 0.24 | U | 6.0 | 0.24 |
| 1634-04-4 | Methyl tert-butyl ether | 0.51 | U | 6.0 | 0.51 |
| 75-09-2 | Methylene Chloride | 0.88 | U | 6.0 | 0.88 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.50 | U | 12 | 0.50 |
| 91-20-3 | Naphthalene | 0.42 | U | 6.0 | 0.42 |
| 104-51-8 | n-Butylbenzene | 0.70 | U | 6.0 | 0.70 |
| 103-65-1 | N-Propylbenzene | 0.71 | U | 6.0 | 0.71 |
| 95-47-6 | o-Xylene | 0.68 | U | 6.0 | 0.68 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 6.0 | 1.1 |
| 100-42-5 | Styrene | 0.48 | U | 6.0 | 0.48 |
| 994-05-8 | Tert-amyl methyl ether | 0.51 | U | 6.0 | 0.51 |
| 637-92-3 | Ethyl t-butyl ether | 0.49 | U | 6.0 | 0.49 |
| 127-18-4 | Tetrachloroethene | 0.95 | U | 6.0 | 0.95 |
| 108-88-3 | Toluene | 0.32 | U | 6.0 | 0.32 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.38 | U | 6.0 | 0.38 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.35 | U | 6.0 | 0.35 |
| 79-01-6 | Trichloroethene | 0.45 | U | 6.0 | 0.45 |
| 75-69-4 | Trichlorofluoromethane | 0.35 | U | 6.0 | 0.35 |
| 108-05-4 | Vinyl acetate | 0.79 | U | 12 | 0.79 |
| 75-01-4 | Vinyl chloride | 0.36 | U | 6.0 | 0.36 |
| 1330-20-7 | Xylenes, Total | 0.64 | U | 12 | 0.64 |
| 98-06-6 | tert-Butylbenzene | 0.80 | U | 6.0 | 0.80 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 146349.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.37(g) | Date Analyzed: 11/10/2015 19:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.72 | U | 5.5 | 0.72 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.5 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.89 | U | 5.5 | 0.89 |
| 563-58-6 | 1,1-Dichloropropene | 0.41 | U | 5.5 | 0.41 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.86 | U | 5.5 | 0.86 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.51 | U | 5.5 | 0.51 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.42 | U | 5.5 | 0.42 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.64 | U | 5.5 | 0.64 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.5 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.51 | U | 5.5 | 0.51 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.76 | U | 5.5 | 0.76 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.5 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 3.7 | J B | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.71 | U | 55 | 0.71 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.75 | U | 22 | 0.75 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.5 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.45 | U | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 6.3 | U | 22 | 6.3 |
| 71-43-2 | Benzene | 1.0 | U | 5.5 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.5 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.5 | 0.27 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.65 | U | 5.5 | 0.65 |
| 56-23-5 | Carbon tetrachloride | 0.72 | U | 5.5 | 0.72 |
| 108-90-7 | Chlorobenzene | 0.59 | U | 5.5 | 0.59 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 146349.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.37(g) | Date Analyzed: 11/10/2015 19:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.41 | U | 5.5 | 0.41 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.5 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.5 | 0.40 |
| 74-87-3 | Chloromethane | 0.87 | U | 5.5 | 0.87 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.79 | U | 5.5 | 0.79 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.5 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.3 | U | 220 | 6.3 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.65 | U | 11 | 0.65 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.5 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.82 | U | 5.5 | 0.82 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.5 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.65 | U | 5.5 | 0.65 |
| 103-65-1 | N-Propylbenzene | 0.66 | U | 5.5 | 0.66 |
| 95-47-6 | o-Xylene | 0.63 | U | 5.5 | 0.63 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.5 | 1.0 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.5 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.89 | U | 5.5 | 0.89 |
| 108-88-3 | Toluene | 0.30 | J | 5.5 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.42 | U | 5.5 | 0.42 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.73 | U | 11 | 0.73 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.60 | U | 11 | 0.60 |
| 98-06-6 | tert-Butylbenzene | 0.74 | U | 5.5 | 0.74 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 146349.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.37(g) | Date Analyzed: 11/10/2015 19:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.59 | U | 5.5 | 0.59 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 99 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 103 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 146350.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 5.67(g) | Date Analyzed: 11/10/2015 19:44

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.52 | U | 5.2 | 0.52 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.67 | U | 5.2 | 0.67 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.32 | U | 5.2 | 0.32 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.43 | U | 5.2 | 0.43 |
| 75-34-3 | 1,1-Dichloroethane | 0.35 | U | 5.2 | 0.35 |
| 75-35-4 | 1,1-Dichloroethene | 0.83 | U | 5.2 | 0.83 |
| 563-58-6 | 1,1-Dichloropropene | 0.38 | U | 5.2 | 0.38 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.52 | U | 5.2 | 0.52 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.81 | U | 5.2 | 0.81 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.2 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.39 | U | 5.2 | 0.39 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.60 | U | 5.2 | 0.60 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.8 | U | 10 | 2.8 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.35 | U | 5.2 | 0.35 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.2 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.2 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.71 | U | 5.2 | 0.71 |
| 142-28-9 | 1,3-Dichloropropane | 0.41 | U | 5.2 | 0.41 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.29 | U | 5.2 | 0.29 |
| 594-20-7 | 2,2-Dichloropropane | 0.45 | U | 5.2 | 0.45 |
| 78-93-3 | 2-Butanone (MEK) | 3.1 | J B | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.66 | U | 52 | 0.66 |
| 95-49-8 | 2-Chlorotoluene | 0.31 | U | 5.2 | 0.31 |
| 591-78-6 | 2-Hexanone | 0.70 | U | 21 | 0.70 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.2 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.42 | U | 5.2 | 0.42 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 5.9 | U | 21 | 5.9 |
| 71-43-2 | Benzene | 0.93 | U | 5.2 | 0.93 |
| 108-86-1 | Bromobenzene | 0.34 | U | 5.2 | 0.34 |
| 75-25-2 | Bromoform | 0.25 | U | 5.2 | 0.25 |
| 74-83-9 | Bromomethane | 0.43 | U | 5.2 | 0.43 |
| 75-15-0 | Carbon disulfide | 0.61 | U | 5.2 | 0.61 |
| 56-23-5 | Carbon tetrachloride | 0.67 | U | 5.2 | 0.67 |
| 108-90-7 | Chlorobenzene | 0.55 | U | 5.2 | 0.55 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 146350.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 5.67(g) | Date Analyzed: 11/10/2015 19:44

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.38 | U | 5.2 | 0.38 |
| 75-00-3 | Chloroethane | 0.45 | U | 5.2 | 0.45 |
| 67-66-3 | Chloroform | 0.37 | U | 5.2 | 0.37 |
| 74-87-3 | Chloromethane | 0.82 | U | 5.2 | 0.82 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.29 | U | 5.2 | 0.29 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.73 | U | 5.2 | 0.73 |
| 74-95-3 | Dibromomethane | 0.43 | U | 5.2 | 0.43 |
| 74-97-5 | Bromochloromethane | 0.34 | U | 5.2 | 0.34 |
| 75-71-8 | Dichlorodifluoromethane | 0.29 | U | 5.2 | 0.29 |
| 75-27-4 | Bromodichloromethane | 0.30 | U | 5.2 | 0.30 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.9 | U | 210 | 5.9 |
| 100-41-4 | Ethylbenzene | 0.33 | U | 5.2 | 0.33 |
| 87-68-3 | Hexachlorobutadiene | 0.41 | U | 5.2 | 0.41 |
| 108-20-3 | Diisopropyl ether | 0.61 | U | 10 | 0.61 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.2 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.44 | U | 5.2 | 0.44 |
| 75-09-2 | Methylene Chloride | 0.76 | U | 5.2 | 0.76 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.43 | U | 10 | 0.43 |
| 91-20-3 | Naphthalene | 0.36 | U | 5.2 | 0.36 |
| 104-51-8 | n-Butylbenzene | 0.61 | U | 5.2 | 0.61 |
| 103-65-1 | N-Propylbenzene | 0.62 | U | 5.2 | 0.62 |
| 95-47-6 | o-Xylene | 0.59 | U | 5.2 | 0.59 |
| 135-98-8 | sec-Butylbenzene | 0.94 | U | 5.2 | 0.94 |
| 100-42-5 | Styrene | 0.41 | U | 5.2 | 0.41 |
| 994-05-8 | Tert-amyl methyl ether | 0.44 | U | 5.2 | 0.44 |
| 637-92-3 | Ethyl t-butyl ether | 0.42 | U | 5.2 | 0.42 |
| 127-18-4 | Tetrachloroethene | 0.83 | U | 5.2 | 0.83 |
| 108-88-3 | Toluene | 0.28 | U | 5.2 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.33 | U | 5.2 | 0.33 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.30 | U | 5.2 | 0.30 |
| 79-01-6 | Trichloroethene | 0.39 | U | 5.2 | 0.39 |
| 75-69-4 | Trichlorofluoromethane | 0.30 | U | 5.2 | 0.30 |
| 108-05-4 | Vinyl acetate | 0.68 | U | 10 | 0.68 |
| 75-01-4 | Vinyl chloride | 0.31 | U | 5.2 | 0.31 |
| 1330-20-7 | Xylenes, Total | 0.56 | U | 10 | 0.56 |
| 98-06-6 | tert-Butylbenzene | 0.69 | U | 5.2 | 0.69 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 146350.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 5.67(g) | Date Analyzed: 11/10/2015 19:44

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.55 | U | 5.2 | 0.55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 105 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 90 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 100 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 146351.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.55(g) | Date Analyzed: 11/10/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.53 | U | 5.3 | 0.53 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.68 | U | 5.3 | 0.68 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.33 | U | 5.3 | 0.33 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.44 | U | 5.3 | 0.44 |
| 75-34-3 | 1,1-Dichloroethane | 0.36 | U | 5.3 | 0.36 |
| 75-35-4 | 1,1-Dichloroethene | 0.84 | U | 5.3 | 0.84 |
| 563-58-6 | 1,1-Dichloropropene | 0.39 | U | 5.3 | 0.39 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.53 | U | 5.3 | 0.53 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.82 | U | 5.3 | 0.82 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.3 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.40 | U | 5.3 | 0.40 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.61 | U | 5.3 | 0.61 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.9 | U | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.36 | U | 5.3 | 0.36 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.3 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.3 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.73 | U | 5.3 | 0.73 |
| 142-28-9 | 1,3-Dichloropropane | 0.42 | U | 5.3 | 0.42 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.30 | U | 5.3 | 0.30 |
| 594-20-7 | 2,2-Dichloropropane | 0.46 | U | 5.3 | 0.46 |
| 78-93-3 | 2-Butanone (MEK) | 4.0 | J B | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.67 | U | 53 | 0.67 |
| 95-49-8 | 2-Chlorotoluene | 0.32 | U | 5.3 | 0.32 |
| 591-78-6 | 2-Hexanone | 0.72 | U | 21 | 0.72 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.3 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 5.3 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 6.0 | U | 21 | 6.0 |
| 71-43-2 | Benzene | 0.95 | U | 5.3 | 0.95 |
| 108-86-1 | Bromobenzene | 0.35 | U | 5.3 | 0.35 |
| 75-25-2 | Bromoform | 0.25 | U | 5.3 | 0.25 |
| 74-83-9 | Bromomethane | 0.44 | U | 5.3 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.62 | U | 5.3 | 0.62 |
| 56-23-5 | Carbon tetrachloride | 0.68 | U | 5.3 | 0.68 |
| 108-90-7 | Chlorobenzene | 0.56 | U | 5.3 | 0.56 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 146351.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.55(g) | Date Analyzed: 11/10/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.39 | U | 5.3 | 0.39 |
| 75-00-3 | Chloroethane | 0.46 | U | 5.3 | 0.46 |
| 67-66-3 | Chloroform | 0.38 | U | 5.3 | 0.38 |
| 74-87-3 | Chloromethane | 0.83 | U | 5.3 | 0.83 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.30 | U | 5.3 | 0.30 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.75 | U | 5.3 | 0.75 |
| 74-95-3 | Dibromomethane | 0.44 | U | 5.3 | 0.44 |
| 74-97-5 | Bromochloromethane | 0.35 | U | 5.3 | 0.35 |
| 75-71-8 | Dichlorodifluoromethane | 0.30 | U | 5.3 | 0.30 |
| 75-27-4 | Bromodichloromethane | 0.31 | U | 5.3 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.0 | U | 210 | 6.0 |
| 100-41-4 | Ethylbenzene | 0.34 | U | 5.3 | 0.34 |
| 87-68-3 | Hexachlorobutadiene | 0.42 | U | 5.3 | 0.42 |
| 108-20-3 | Diisopropyl ether | 0.62 | U | 11 | 0.62 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.3 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.45 | U | 5.3 | 0.45 |
| 75-09-2 | Methylene Chloride | 0.78 | U | 5.3 | 0.78 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.44 | U | 11 | 0.44 |
| 91-20-3 | Naphthalene | 0.37 | U | 5.3 | 0.37 |
| 104-51-8 | n-Butylbenzene | 0.62 | U | 5.3 | 0.62 |
| 103-65-1 | N-Propylbenzene | 0.63 | U | 5.3 | 0.63 |
| 95-47-6 | o-Xylene | 0.60 | U | 5.3 | 0.60 |
| 135-98-8 | sec-Butylbenzene | 0.96 | U | 5.3 | 0.96 |
| 100-42-5 | Styrene | 0.42 | U | 5.3 | 0.42 |
| 994-05-8 | Tert-amyl methyl ether | 0.45 | U | 5.3 | 0.45 |
| 637-92-3 | Ethyl t-butyl ether | 0.43 | U | 5.3 | 0.43 |
| 127-18-4 | Tetrachloroethene | 0.84 | U | 5.3 | 0.84 |
| 108-88-3 | Toluene | 0.29 | J | 5.3 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.34 | U | 5.3 | 0.34 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.31 | U | 5.3 | 0.31 |
| 79-01-6 | Trichloroethene | 0.40 | U | 5.3 | 0.40 |
| 75-69-4 | Trichlorofluoromethane | 0.31 | U | 5.3 | 0.31 |
| 108-05-4 | Vinyl acetate | 0.70 | U | 11 | 0.70 |
| 75-01-4 | Vinyl chloride | 0.32 | U | 5.3 | 0.32 |
| 1330-20-7 | Xylenes, Total | 0.57 | U | 11 | 0.57 |
| 98-06-6 | tert-Butylbenzene | 0.71 | U | 5.3 | 0.71 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 146351.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.55(g) | Date Analyzed: 11/10/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.56 | U | 5.3 | 0.56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 87 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 89 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 78 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 96 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 146365.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.55(g) | Date Analyzed: 11/11/2015 13:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.53 | U | 5.3 | 0.53 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.68 | U | 5.3 | 0.68 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.33 | U | 5.3 | 0.33 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.44 | U | 5.3 | 0.44 |
| 75-34-3 | 1,1-Dichloroethane | 0.36 | U | 5.3 | 0.36 |
| 75-35-4 | 1,1-Dichloroethene | 0.84 | U | 5.3 | 0.84 |
| 563-58-6 | 1,1-Dichloropropene | 0.39 | U | 5.3 | 0.39 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.53 | U | 5.3 | 0.53 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.82 | U | 5.3 | 0.82 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.3 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.40 | U | 5.3 | 0.40 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.61 | U | 5.3 | 0.61 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.9 | U | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.36 | U | 5.3 | 0.36 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.3 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.3 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.73 | U | 5.3 | 0.73 |
| 142-28-9 | 1,3-Dichloropropane | 0.42 | U | 5.3 | 0.42 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.29 | U | 5.3 | 0.29 |
| 594-20-7 | 2,2-Dichloropropane | 0.46 | U | 5.3 | 0.46 |
| 78-93-3 | 2-Butanone (MEK) | 3.6 | J B | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.67 | U | 53 | 0.67 |
| 95-49-8 | 2-Chlorotoluene | 0.32 | U | 5.3 | 0.32 |
| 591-78-6 | 2-Hexanone | 0.72 | U | 21 | 0.72 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.3 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 5.3 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 6.0 | U | 21 | 6.0 |
| 71-43-2 | Benzene | 0.95 | U | 5.3 | 0.95 |
| 108-86-1 | Bromobenzene | 0.35 | U | 5.3 | 0.35 |
| 75-25-2 | Bromoform | 0.25 | U | 5.3 | 0.25 |
| 74-83-9 | Bromomethane | 0.44 | U | 5.3 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.62 | U | 5.3 | 0.62 |
| 56-23-5 | Carbon tetrachloride | 0.68 | U | 5.3 | 0.68 |
| 108-90-7 | Chlorobenzene | 0.56 | U | 5.3 | 0.56 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 146365.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.55(g) | Date Analyzed: 11/11/2015 13:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.39 | U | 5.3 | 0.39 |
| 75-00-3 | Chloroethane | 0.46 | U | 5.3 | 0.46 |
| 67-66-3 | Chloroform | 0.38 | U | 5.3 | 0.38 |
| 74-87-3 | Chloromethane | 0.83 | U | 5.3 | 0.83 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.29 | U | 5.3 | 0.29 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.75 | U | 5.3 | 0.75 |
| 74-95-3 | Dibromomethane | 0.44 | U | 5.3 | 0.44 |
| 74-97-5 | Bromochloromethane | 0.35 | U | 5.3 | 0.35 |
| 75-71-8 | Dichlorodifluoromethane | 0.29 | U | 5.3 | 0.29 |
| 75-27-4 | Bromodichloromethane | 0.31 | U | 5.3 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.0 | U | 210 | 6.0 |
| 100-41-4 | Ethylbenzene | 0.34 | U | 5.3 | 0.34 |
| 87-68-3 | Hexachlorobutadiene | 0.42 | U | 5.3 | 0.42 |
| 108-20-3 | Diisopropyl ether | 0.62 | U | 11 | 0.62 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.3 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.45 | U | 5.3 | 0.45 |
| 75-09-2 | Methylene Chloride | 0.78 | U | 5.3 | 0.78 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.44 | U | 11 | 0.44 |
| 91-20-3 | Naphthalene | 0.37 | U | 5.3 | 0.37 |
| 104-51-8 | n-Butylbenzene | 0.62 | U | 5.3 | 0.62 |
| 103-65-1 | N-Propylbenzene | 0.63 | U | 5.3 | 0.63 |
| 95-47-6 | o-Xylene | 0.60 | U | 5.3 | 0.60 |
| 135-98-8 | sec-Butylbenzene | 0.96 | U | 5.3 | 0.96 |
| 100-42-5 | Styrene | 0.42 | U | 5.3 | 0.42 |
| 994-05-8 | Tert-amyl methyl ether | 0.45 | U | 5.3 | 0.45 |
| 637-92-3 | Ethyl t-butyl ether | 0.43 | U | 5.3 | 0.43 |
| 127-18-4 | Tetrachloroethene | 0.84 | U | 5.3 | 0.84 |
| 108-88-3 | Toluene | 0.28 | U | 5.3 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.34 | U | 5.3 | 0.34 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.31 | U | 5.3 | 0.31 |
| 79-01-6 | Trichloroethene | 0.40 | U | 5.3 | 0.40 |
| 75-69-4 | Trichlorofluoromethane | 0.31 | U | 5.3 | 0.31 |
| 108-05-4 | Vinyl acetate | 0.70 | U | 11 | 0.70 |
| 75-01-4 | Vinyl chloride | 0.32 | U | 5.3 | 0.32 |
| 1330-20-7 | Xylenes, Total | 0.57 | U | 11 | 0.57 |
| 98-06-6 | tert-Butylbenzene | 0.71 | U | 5.3 | 0.71 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 146365.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.55(g) | Date Analyzed: 11/11/2015 13:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.56 | U | 5.3 | 0.56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 102 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 89 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: TB-110515 | Lab Sample ID: 240-57597-16

Matrix: Water | Lab File ID: UXC4397.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.28 | U | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.44 | U | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.22 | U | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.30 | U | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 0.45 | U | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.37 | U | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.44 | U | 1.0 | 0.44 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.47 | U | 5.0 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.32 | U | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.41 | U | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.82 | U | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.25 | U | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 0.23 | U | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 0.25 | U | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.19 | U | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 0.19 | U | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.27 | U | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 0.26 | U | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 0.53 | U | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.59 | U | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 0.40 | U | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 0.48 | U | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 10 | 0.99 |
| 67-64-1 | Acetone | 0.94 | U | 10 | 0.94 |
| 71-43-2 | Benzene | 0.35 | U | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 0.35 | U | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 0.56 | U | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 0.44 | U | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.38 | U | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 0.43 | U | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 0.25 | U | 1.0 | 0.25 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: TB-110515 | Lab Sample ID: 240-57597-16

Matrix: Water | Lab File ID: UXC4397.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 1.0 | 0.43 |
| 75-00-3 | Chloroethane | 0.32 | U | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 0.25 | U | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 0.44 | U | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.26 | U | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.46 | U | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 0.42 | U | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.50 | U | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 4.9 | U | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 0.25 | U | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 0.35 | U | 1.0 | 0.35 |
| 108-20-3 | Diisopropyl ether | 0.50 | U | 10 | 0.50 |
| 98-82-8 | Isopropylbenzene | 0.35 | U | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 0.20 | U | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 0.33 | U | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.24 | U | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 0.45 | U | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 0.31 | U | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 0.40 | U | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 0.25 | U | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 0.48 | U | 1.0 | 0.48 |
| 100-42-5 | Styrene | 0.45 | U | 1.0 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.30 | U | 5.0 | 0.30 |
| 637-92-3 | Ethyl t-butyl ether | 0.23 | U | 5.0 | 0.23 |
| 127-18-4 | Tetrachloroethene | 0.31 | U | 1.0 | 0.31 |
| 108-88-3 | Toluene | 0.23 | U | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.30 | U | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.56 | U | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 0.22 | U | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 0.49 | U | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 0.41 | U | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 0.29 | U | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 0.52 | U | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 0.41 | U | 1.0 | 0.41 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: TB-110515 | Lab Sample ID: 240-57597-16

Matrix: Water | Lab File ID: UXC4397.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.32 | U | 1.0 | 0.32 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 85 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 97 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 90 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 99 | | 78-125 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 146336.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 5.89(g) | Date Analyzed: 11/10/2015 14:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.81 | U | 5.0 | 0.81 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.79 | U | 5.0 | 0.79 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.91 | U | 5.0 | 0.91 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 146336.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 5.89(g) | Date Analyzed: 11/10/2015 14:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.80 | U | 5.0 | 0.80 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.72 | U | 5.0 | 0.72 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.8 | U | 200 | 5.8 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.75 | U | 5.0 | 0.75 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.35 | U | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.92 | U | 5.0 | 0.92 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.81 | U | 5.0 | 0.81 |
| 108-88-3 | Toluene | 0.27 | U | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: 146336.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 5.89(g) | Date Analyzed: 11/10/2015 14:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 103 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 91 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 103 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 146337.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 5.36(g) | Date Analyzed: 11/10/2015 15:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.71 | U | 5.5 | 0.71 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.37 | U | 5.5 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 0.88 | U | 5.5 | 0.88 |
| 563-58-6 | 1,1-Dichloropropene | 0.41 | U | 5.5 | 0.41 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.86 | U | 5.5 | 0.86 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.50 | U | 5.5 | 0.50 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.42 | U | 5.5 | 0.42 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.64 | U | 5.5 | 0.64 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.37 | U | 5.5 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 0.50 | U | 5.5 | 0.50 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.76 | U | 5.5 | 0.76 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.48 | U | 5.5 | 0.48 |
| 78-93-3 | 2-Butanone (MEK) | 4.0 | J B | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.70 | U | 55 | 0.70 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.75 | U | 22 | 0.75 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.5 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.45 | U | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 22 | 1.3 |
| 67-64-1 | Acetone | 6.3 | U | 22 | 6.3 |
| 71-43-2 | Benzene | 0.99 | U | 5.5 | 0.99 |
| 108-86-1 | Bromobenzene | 0.36 | U | 5.5 | 0.36 |
| 75-25-2 | Bromoform | 0.26 | U | 5.5 | 0.26 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.65 | U | 5.5 | 0.65 |
| 56-23-5 | Carbon tetrachloride | 0.71 | U | 5.5 | 0.71 |
| 108-90-7 | Chlorobenzene | 0.58 | U | 5.5 | 0.58 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 146337.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 5.36(g) | Date Analyzed: 11/10/2015 15:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.41 | U | 5.5 | 0.41 |
| 75-00-3 | Chloroethane | 0.48 | U | 5.5 | 0.48 |
| 67-66-3 | Chloroform | 0.39 | U | 5.5 | 0.39 |
| 74-87-3 | Chloromethane | 0.87 | U | 5.5 | 0.87 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.78 | U | 5.5 | 0.78 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.36 | U | 5.5 | 0.36 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.3 | U | 220 | 6.3 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.65 | U | 11 | 0.65 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.47 | U | 5.5 | 0.47 |
| 75-09-2 | Methylene Chloride | 0.81 | U | 5.5 | 0.81 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.38 | U | 5.5 | 0.38 |
| 104-51-8 | n-Butylbenzene | 0.65 | U | 5.5 | 0.65 |
| 103-65-1 | N-Propylbenzene | 0.66 | U | 5.5 | 0.66 |
| 95-47-6 | o-Xylene | 0.63 | U | 5.5 | 0.63 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.5 | 1.0 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.47 | U | 5.5 | 0.47 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.88 | U | 5.5 | 0.88 |
| 108-88-3 | Toluene | 0.30 | U | 5.5 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.42 | U | 5.5 | 0.42 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.72 | U | 11 | 0.72 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.59 | U | 11 | 0.59 |
| 98-06-6 | tert-Butylbenzene | 0.74 | U | 5.5 | 0.74 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: 146337.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 5.36(g) | Date Analyzed: 11/10/2015 15:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.58 | U | 5.5 | 0.58 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 96 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 95 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 85 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 99 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 146338.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 5.23(g) | Date Analyzed: 11/10/2015 15:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.76 | U | 5.8 | 0.76 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 5.8 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.91 | U | 5.8 | 0.91 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 5.8 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 5.8 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.48 | U | 5.8 | 0.48 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 6.6 | U | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.69 | U | 5.8 | 0.69 |
| 56-23-5 | Carbon tetrachloride | 0.76 | U | 5.8 | 0.76 |
| 108-90-7 | Chlorobenzene | 0.62 | U | 5.8 | 0.62 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 146338.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 5.23(g) | Date Analyzed: 11/10/2015 15:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 5.8 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.83 | U | 5.8 | 0.83 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 5.8 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.69 | U | 12 | 0.69 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.41 | U | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 0.69 | U | 5.8 | 0.69 |
| 103-65-1 | N-Propylbenzene | 0.70 | U | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.48 | U | 5.8 | 0.48 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.77 | U | 12 | 0.77 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.63 | U | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 0.78 | U | 5.8 | 0.78 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: 146338.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 5.23(g) | Date Analyzed: 11/10/2015 15:26

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.62 | U | 5.8 | 0.62 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 93 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 90 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 86 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 90 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 146339.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 5.43(g) | Date Analyzed: 11/10/2015 15:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.71 | U | 5.5 | 0.71 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.37 | U | 5.5 | 0.37 |
| 75-35-4 | 1,1-Dichloroethene | 0.87 | U | 5.5 | 0.87 |
| 563-58-6 | 1,1-Dichloropropene | 0.40 | U | 5.5 | 0.40 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.85 | U | 5.5 | 0.85 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.50 | U | 5.5 | 0.50 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.41 | U | 5.5 | 0.41 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.63 | U | 5.5 | 0.63 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.37 | U | 5.5 | 0.37 |
| 107-06-2 | 1,2-Dichloroethane | 0.50 | U | 5.5 | 0.50 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 5.5 | 0.75 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.48 | U | 5.5 | 0.48 |
| 78-93-3 | 2-Butanone (MEK) | 4.0 | J B | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.70 | U | 55 | 0.70 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.74 | U | 22 | 0.74 |
| 106-43-4 | 4-Chlorotoluene | 0.28 | U | 5.5 | 0.28 |
| 99-87-6 | p-Isopropyltoluene | 0.45 | U | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 22 | 1.3 |
| 67-64-1 | Acetone | 6.2 | U | 22 | 6.2 |
| 71-43-2 | Benzene | 0.98 | U | 5.5 | 0.98 |
| 108-86-1 | Bromobenzene | 0.36 | U | 5.5 | 0.36 |
| 75-25-2 | Bromoform | 0.26 | U | 5.5 | 0.26 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.64 | U | 5.5 | 0.64 |
| 56-23-5 | Carbon tetrachloride | 0.71 | U | 5.5 | 0.71 |
| 108-90-7 | Chlorobenzene | 0.58 | U | 5.5 | 0.58 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 146339.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 5.43(g) | Date Analyzed: 11/10/2015 15:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.40 | U | 5.5 | 0.40 |
| 75-00-3 | Chloroethane | 0.48 | U | 5.5 | 0.48 |
| 67-66-3 | Chloroform | 0.39 | U | 5.5 | 0.39 |
| 74-87-3 | Chloromethane | 0.86 | U | 5.5 | 0.86 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.77 | U | 5.5 | 0.77 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.36 | U | 5.5 | 0.36 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.2 | U | 220 | 6.2 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.64 | U | 11 | 0.64 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.47 | U | 5.5 | 0.47 |
| 75-09-2 | Methylene Chloride | 0.81 | U | 5.5 | 0.81 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.38 | U | 5.5 | 0.38 |
| 104-51-8 | n-Butylbenzene | 0.64 | U | 5.5 | 0.64 |
| 103-65-1 | N-Propylbenzene | 0.65 | U | 5.5 | 0.65 |
| 95-47-6 | o-Xylene | 0.62 | U | 5.5 | 0.62 |
| 135-98-8 | sec-Butylbenzene | 0.99 | U | 5.5 | 0.99 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.47 | U | 5.5 | 0.47 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.87 | U | 5.5 | 0.87 |
| 108-88-3 | Toluene | 0.29 | U | 5.5 | 0.29 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.41 | U | 5.5 | 0.41 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.72 | U | 11 | 0.72 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.59 | U | 11 | 0.59 |
| 98-06-6 | tert-Butylbenzene | 0.73 | U | 5.5 | 0.73 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: 146339.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 5.43(g) | Date Analyzed: 11/10/2015 15:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.58 | U | 5.5 | 0.58 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 93 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 94 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 83 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 146340.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.09(g) | Date Analyzed: 11/10/2015 16:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U F2 | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.3 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U F2 | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U F1 | 23 | 1.4 |
| 67-64-1 | Acetone | 13 | J | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 146340.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.09(g) | Date Analyzed: 11/10/2015 16:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.92 | U | 5.8 | 0.92 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U F2 | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.41 | J | 5.8 | 0.41 |
| 104-51-8 | n-Butylbenzene | 0.68 | U F2 | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.70 | U | 5.8 | 0.70 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U F2 | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U F1 | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.63 | U | 12 | 0.63 |
| 98-06-6 | tert-Butylbenzene | 0.78 | U | 5.8 | 0.78 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: 146340.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 5.09(g) | Date Analyzed: 11/10/2015 16:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 100 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 104 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 146343.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.18(g) | Date Analyzed: 11/10/2015 17:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.93 | U | 5.8 | 0.93 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 3.8 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 6.6 | U | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 146343.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.18(g) | Date Analyzed: 11/10/2015 17:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.86 | U | 5.8 | 0.86 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.45 | J | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.93 | U | 5.8 | 0.93 |
| 108-88-3 | Toluene | 0.31 | U | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: 146343.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.18(g) | Date Analyzed: 11/10/2015 17:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 88 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 89 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 81 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 146344.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 5.31(g) | Date Analyzed: 11/10/2015 17:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.56 | U | 5.6 | 0.56 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.73 | U | 5.6 | 0.73 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.35 | U | 5.6 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.47 | U | 5.6 | 0.47 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.6 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.90 | U | 5.6 | 0.90 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 5.6 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.56 | U | 5.6 | 0.56 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.88 | U | 5.6 | 0.88 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.52 | U | 5.6 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.43 | U | 5.6 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.65 | U | 5.6 | 0.65 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.6 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.52 | U | 5.6 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.6 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.78 | U | 5.6 | 0.78 |
| 142-28-9 | 1,3-Dichloropropane | 0.45 | U | 5.6 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.6 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.6 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 4.1 | J B | 22 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.72 | U | 56 | 0.72 |
| 95-49-8 | 2-Chlorotoluene | 0.34 | U | 5.6 | 0.34 |
| 591-78-6 | 2-Hexanone | 0.76 | U | 22 | 0.76 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.6 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.46 | U | 5.6 | 0.46 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 6.4 | U | 22 | 6.4 |
| 71-43-2 | Benzene | 1.0 | U | 5.6 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.6 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.6 | 0.27 |
| 74-83-9 | Bromomethane | 0.47 | U | 5.6 | 0.47 |
| 75-15-0 | Carbon disulfide | 0.66 | U | 5.6 | 0.66 |
| 56-23-5 | Carbon tetrachloride | 0.73 | U | 5.6 | 0.73 |
| 108-90-7 | Chlorobenzene | 0.60 | U | 5.6 | 0.60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 146344.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 5.31(g) | Date Analyzed: 11/10/2015 17:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.42 | U | 5.6 | 0.42 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.6 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.6 | 0.40 |
| 74-87-3 | Chloromethane | 0.89 | U | 5.6 | 0.89 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.6 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.80 | U | 5.6 | 0.80 |
| 74-95-3 | Dibromomethane | 0.47 | U | 5.6 | 0.47 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.6 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.6 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.6 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.4 | U | 220 | 6.4 |
| 100-41-4 | Ethylbenzene | 0.36 | U | 5.6 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 0.45 | U | 5.6 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.66 | U | 11 | 0.66 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.6 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.6 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.83 | U | 5.6 | 0.83 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.47 | U | 11 | 0.47 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.6 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.66 | U | 5.6 | 0.66 |
| 103-65-1 | N-Propylbenzene | 0.67 | U | 5.6 | 0.67 |
| 95-47-6 | o-Xylene | 0.64 | U | 5.6 | 0.64 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.6 | 1.0 |
| 100-42-5 | Styrene | 0.45 | U | 5.6 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.6 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.46 | U | 5.6 | 0.46 |
| 127-18-4 | Tetrachloroethene | 0.90 | U | 5.6 | 0.90 |
| 108-88-3 | Toluene | 0.30 | U | 5.6 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.36 | U | 5.6 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.6 | 0.33 |
| 79-01-6 | Trichloroethene | 0.43 | U | 5.6 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.6 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.74 | U | 11 | 0.74 |
| 75-01-4 | Vinyl chloride | 0.34 | U | 5.6 | 0.34 |
| 1330-20-7 | Xylenes, Total | 0.61 | U | 11 | 0.61 |
| 98-06-6 | tert-Butylbenzene | 0.75 | U | 5.6 | 0.75 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: 146344.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:12

Sample wt/vol: 5.31(g) | Date Analyzed: 11/10/2015 17:35

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.2 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.60 | U | 5.6 | 0.60 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 98 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 103 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 146345.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 5.29(g) | Date Analyzed: 11/10/2015 17:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.48 | U | 5.8 | 0.48 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.92 | U | 5.8 | 0.92 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.3 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.78 | U | 23 | 0.78 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 6.6 | U | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.48 | U | 5.8 | 0.48 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 146345.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 5.29(g) | Date Analyzed: 11/10/2015 17:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.41 | U | 5.8 | 0.41 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.48 | U | 5.8 | 0.48 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.8 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.85 | U | 5.8 | 0.85 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.48 | U | 12 | 0.48 |
| 91-20-3 | Naphthalene | 0.40 | U | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.8 | 1.0 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.92 | U | 5.8 | 0.92 |
| 108-88-3 | Toluene | 0.33 | J | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.8 | 0.33 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.8 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: 146345.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 5.29(g) | Date Analyzed: 11/10/2015 17:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 97 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 93 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 85 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 146346.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.35(g) | Date Analyzed: 11/10/2015 18:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.58 | U | 5.8 | 0.58 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.75 | U | 5.8 | 0.75 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.36 | U | 5.8 | 0.36 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.49 | U | 5.8 | 0.49 |
| 75-34-3 | 1,1-Dichloroethane | 0.39 | U | 5.8 | 0.39 |
| 75-35-4 | 1,1-Dichloroethene | 0.92 | U | 5.8 | 0.92 |
| 563-58-6 | 1,1-Dichloropropene | 0.43 | U | 5.8 | 0.43 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.58 | U | 5.8 | 0.58 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.90 | U | 5.8 | 0.90 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.53 | U | 5.8 | 0.53 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.44 | U | 5.8 | 0.44 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.67 | U | 5.8 | 0.67 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.1 | U | 12 | 3.1 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.39 | U | 5.8 | 0.39 |
| 107-06-2 | 1,2-Dichloroethane | 0.53 | U | 5.8 | 0.53 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.8 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 5.8 | 0.80 |
| 142-28-9 | 1,3-Dichloropropane | 0.46 | U | 5.8 | 0.46 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.32 | U | 5.8 | 0.32 |
| 594-20-7 | 2,2-Dichloropropane | 0.51 | U | 5.8 | 0.51 |
| 78-93-3 | 2-Butanone (MEK) | 4.4 | J B | 23 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.74 | U | 58 | 0.74 |
| 95-49-8 | 2-Chlorotoluene | 0.35 | U | 5.8 | 0.35 |
| 591-78-6 | 2-Hexanone | 0.79 | U | 23 | 0.79 |
| 106-43-4 | 4-Chlorotoluene | 0.30 | U | 5.8 | 0.30 |
| 99-87-6 | p-Isopropyltoluene | 0.47 | U | 5.8 | 0.47 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 23 | 1.4 |
| 67-64-1 | Acetone | 12 | J | 23 | 6.6 |
| 71-43-2 | Benzene | 1.0 | U | 5.8 | 1.0 |
| 108-86-1 | Bromobenzene | 0.38 | U | 5.8 | 0.38 |
| 75-25-2 | Bromoform | 0.28 | U | 5.8 | 0.28 |
| 74-83-9 | Bromomethane | 0.49 | U | 5.8 | 0.49 |
| 75-15-0 | Carbon disulfide | 0.68 | U | 5.8 | 0.68 |
| 56-23-5 | Carbon tetrachloride | 0.75 | U | 5.8 | 0.75 |
| 108-90-7 | Chlorobenzene | 0.61 | U | 5.8 | 0.61 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 146346.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.35(g) | Date Analyzed: 11/10/2015 18:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 5.8 | 0.43 |
| 75-00-3 | Chloroethane | 0.51 | U | 5.8 | 0.51 |
| 67-66-3 | Chloroform | 0.42 | U | 5.8 | 0.42 |
| 74-87-3 | Chloromethane | 0.91 | U | 5.8 | 0.91 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.32 | U | 5.8 | 0.32 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.82 | U | 5.8 | 0.82 |
| 74-95-3 | Dibromomethane | 0.49 | U | 5.8 | 0.49 |
| 74-97-5 | Bromochloromethane | 0.38 | U | 5.8 | 0.38 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 5.8 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.34 | U | 5.8 | 0.34 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.6 | U | 230 | 6.6 |
| 100-41-4 | Ethylbenzene | 0.37 | U | 5.8 | 0.37 |
| 87-68-3 | Hexachlorobutadiene | 0.46 | U | 5.8 | 0.46 |
| 108-20-3 | Diisopropyl ether | 0.68 | U | 12 | 0.68 |
| 98-82-8 | Isopropylbenzene | 0.23 | U | 5.8 | 0.23 |
| 1634-04-4 | Methyl tert-butyl ether | 0.50 | U | 5.8 | 0.50 |
| 75-09-2 | Methylene Chloride | 0.85 | U | 5.8 | 0.85 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.49 | U | 12 | 0.49 |
| 91-20-3 | Naphthalene | 0.40 | U | 5.8 | 0.40 |
| 104-51-8 | n-Butylbenzene | 0.68 | U | 5.8 | 0.68 |
| 103-65-1 | N-Propylbenzene | 0.69 | U | 5.8 | 0.69 |
| 95-47-6 | o-Xylene | 0.66 | U | 5.8 | 0.66 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 5.8 | 1.1 |
| 100-42-5 | Styrene | 0.46 | U | 5.8 | 0.46 |
| 994-05-8 | Tert-amyl methyl ether | 0.50 | U | 5.8 | 0.50 |
| 637-92-3 | Ethyl t-butyl ether | 0.47 | U | 5.8 | 0.47 |
| 127-18-4 | Tetrachloroethene | 0.92 | U | 5.8 | 0.92 |
| 108-88-3 | Toluene | 0.36 | J | 5.8 | 0.31 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.37 | U | 5.8 | 0.37 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.34 | U | 5.8 | 0.34 |
| 79-01-6 | Trichloroethene | 0.44 | U | 5.8 | 0.44 |
| 75-69-4 | Trichlorofluoromethane | 0.34 | U | 5.8 | 0.34 |
| 108-05-4 | Vinyl acetate | 0.76 | U | 12 | 0.76 |
| 75-01-4 | Vinyl chloride | 0.35 | U | 5.8 | 0.35 |
| 1330-20-7 | Xylenes, Total | 0.62 | U | 12 | 0.62 |
| 98-06-6 | tert-Butylbenzene | 0.77 | U | 5.8 | 0.77 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: 146346.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.35(g) | Date Analyzed: 11/10/2015 18:18

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.61 | U | 5.8 | 0.61 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 88 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 82 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 94 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 146347.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.14(g) | Date Analyzed: 11/10/2015 18:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.56 | U | 5.6 | 0.56 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.73 | U | 5.6 | 0.73 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.35 | U | 5.6 | 0.35 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.47 | U | 5.6 | 0.47 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.6 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.90 | U | 5.6 | 0.90 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 5.6 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.56 | U | 5.6 | 0.56 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.88 | U | 5.6 | 0.88 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.52 | U | 5.6 | 0.52 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.43 | U | 5.6 | 0.43 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.65 | U | 5.6 | 0.65 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.6 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.52 | U | 5.6 | 0.52 |
| 78-87-5 | 1,2-Dichloropropane | 0.16 | U | 5.6 | 0.16 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.78 | U | 5.6 | 0.78 |
| 142-28-9 | 1,3-Dichloropropane | 0.45 | U | 5.6 | 0.45 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.6 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.6 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 3.5 | J B | 22 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.72 | U | 56 | 0.72 |
| 95-49-8 | 2-Chlorotoluene | 0.34 | U | 5.6 | 0.34 |
| 591-78-6 | 2-Hexanone | 0.76 | U | 22 | 0.76 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.6 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.46 | U | 5.6 | 0.46 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 6.4 | U | 22 | 6.4 |
| 71-43-2 | Benzene | 1.0 | U | 5.6 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.6 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.6 | 0.27 |
| 74-83-9 | Bromomethane | 0.47 | U | 5.6 | 0.47 |
| 75-15-0 | Carbon disulfide | 0.66 | U | 5.6 | 0.66 |
| 56-23-5 | Carbon tetrachloride | 0.73 | U | 5.6 | 0.73 |
| 108-90-7 | Chlorobenzene | 0.60 | U | 5.6 | 0.60 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 146347.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.14(g) | Date Analyzed: 11/10/2015 18:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.42 | U | 5.6 | 0.42 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.6 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.6 | 0.40 |
| 74-87-3 | Chloromethane | 0.89 | U | 5.6 | 0.89 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.6 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.80 | U | 5.6 | 0.80 |
| 74-95-3 | Dibromomethane | 0.47 | U | 5.6 | 0.47 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.6 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.6 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.33 | U | 5.6 | 0.33 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.4 | U | 220 | 6.4 |
| 100-41-4 | Ethylbenzene | 0.36 | U | 5.6 | 0.36 |
| 87-68-3 | Hexachlorobutadiene | 0.45 | U | 5.6 | 0.45 |
| 108-20-3 | Diisopropyl ether | 0.66 | U | 11 | 0.66 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.6 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.6 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.83 | U | 5.6 | 0.83 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.47 | U | 11 | 0.47 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.6 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.66 | U | 5.6 | 0.66 |
| 103-65-1 | N-Propylbenzene | 0.67 | U | 5.6 | 0.67 |
| 95-47-6 | o-Xylene | 0.64 | U | 5.6 | 0.64 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.6 | 1.0 |
| 100-42-5 | Styrene | 0.45 | U | 5.6 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.6 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.46 | U | 5.6 | 0.46 |
| 127-18-4 | Tetrachloroethene | 0.90 | U | 5.6 | 0.90 |
| 108-88-3 | Toluene | 0.30 | U | 5.6 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.36 | U | 5.6 | 0.36 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.33 | U | 5.6 | 0.33 |
| 79-01-6 | Trichloroethene | 0.43 | U | 5.6 | 0.43 |
| 75-69-4 | Trichlorofluoromethane | 0.33 | U | 5.6 | 0.33 |
| 108-05-4 | Vinyl acetate | 0.74 | U | 11 | 0.74 |
| 75-01-4 | Vinyl chloride | 0.34 | U | 5.6 | 0.34 |
| 1330-20-7 | Xylenes, Total | 0.61 | U | 11 | 0.61 |
| 98-06-6 | tert-Butylbenzene | 0.75 | U | 5.6 | 0.75 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: 146347.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.14(g) | Date Analyzed: 11/10/2015 18:39

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.60 | U | 5.6 | 0.60 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 89 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 88 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 83 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 93 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 146348.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/10/2015 19:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.60 | U | 6.0 | 0.60 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.77 | U | 6.0 | 0.77 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.37 | U | 6.0 | 0.37 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.50 | U | 6.0 | 0.50 |
| 75-34-3 | 1,1-Dichloroethane | 0.40 | U | 6.0 | 0.40 |
| 75-35-4 | 1,1-Dichloroethene | 0.95 | U | 6.0 | 0.95 |
| 563-58-6 | 1,1-Dichloropropene | 0.44 | U | 6.0 | 0.44 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.60 | U | 6.0 | 0.60 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.93 | U | 6.0 | 0.93 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.55 | U | 6.0 | 0.55 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.45 | U | 6.0 | 0.45 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.69 | U | 6.0 | 0.69 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.2 | U | 12 | 3.2 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.40 | U | 6.0 | 0.40 |
| 107-06-2 | 1,2-Dichloroethane | 0.55 | U | 6.0 | 0.55 |
| 78-87-5 | 1,2-Dichloropropane | 0.17 | U | 6.0 | 0.17 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.82 | U | 6.0 | 0.82 |
| 142-28-9 | 1,3-Dichloropropane | 0.48 | U | 6.0 | 0.48 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.33 | U | 6.0 | 0.33 |
| 594-20-7 | 2,2-Dichloropropane | 0.52 | U | 6.0 | 0.52 |
| 78-93-3 | 2-Butanone (MEK) | 3.9 | J B | 24 | 1.3 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.76 | U | 60 | 0.76 |
| 95-49-8 | 2-Chlorotoluene | 0.36 | U | 6.0 | 0.36 |
| 591-78-6 | 2-Hexanone | 0.81 | U | 24 | 0.81 |
| 106-43-4 | 4-Chlorotoluene | 0.31 | U | 6.0 | 0.31 |
| 99-87-6 | p-Isopropyltoluene | 0.49 | U | 6.0 | 0.49 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.5 | U | 24 | 1.5 |
| 67-64-1 | Acetone | 6.8 | U | 24 | 6.8 |
| 71-43-2 | Benzene | 1.1 | U | 6.0 | 1.1 |
| 108-86-1 | Bromobenzene | 0.39 | U | 6.0 | 0.39 |
| 75-25-2 | Bromoform | 0.29 | U | 6.0 | 0.29 |
| 74-83-9 | Bromomethane | 0.50 | U | 6.0 | 0.50 |
| 75-15-0 | Carbon disulfide | 0.70 | U | 6.0 | 0.70 |
| 56-23-5 | Carbon tetrachloride | 0.77 | U | 6.0 | 0.77 |
| 108-90-7 | Chlorobenzene | 0.63 | U | 6.0 | 0.63 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 146348.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/10/2015 19:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.44 | U | 6.0 | 0.44 |
| 75-00-3 | Chloroethane | 0.52 | U | 6.0 | 0.52 |
| 67-66-3 | Chloroform | 0.43 | U | 6.0 | 0.43 |
| 74-87-3 | Chloromethane | 0.94 | U | 6.0 | 0.94 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.33 | U | 6.0 | 0.33 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.85 | U | 6.0 | 0.85 |
| 74-95-3 | Dibromomethane | 0.50 | U | 6.0 | 0.50 |
| 74-97-5 | Bromochloromethane | 0.39 | U | 6.0 | 0.39 |
| 75-71-8 | Dichlorodifluoromethane | 0.33 | U | 6.0 | 0.33 |
| 75-27-4 | Bromodichloromethane | 0.35 | U | 6.0 | 0.35 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.8 | U | 240 | 6.8 |
| 100-41-4 | Ethylbenzene | 0.38 | U | 6.0 | 0.38 |
| 87-68-3 | Hexachlorobutadiene | 0.48 | U | 6.0 | 0.48 |
| 108-20-3 | Diisopropyl ether | 0.70 | U | 12 | 0.70 |
| 98-82-8 | Isopropylbenzene | 0.24 | U | 6.0 | 0.24 |
| 1634-04-4 | Methyl tert-butyl ether | 0.51 | U | 6.0 | 0.51 |
| 75-09-2 | Methylene Chloride | 0.88 | U | 6.0 | 0.88 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.50 | U | 12 | 0.50 |
| 91-20-3 | Naphthalene | 0.42 | U | 6.0 | 0.42 |
| 104-51-8 | n-Butylbenzene | 0.70 | U | 6.0 | 0.70 |
| 103-65-1 | N-Propylbenzene | 0.71 | U | 6.0 | 0.71 |
| 95-47-6 | o-Xylene | 0.68 | U | 6.0 | 0.68 |
| 135-98-8 | sec-Butylbenzene | 1.1 | U | 6.0 | 1.1 |
| 100-42-5 | Styrene | 0.48 | U | 6.0 | 0.48 |
| 994-05-8 | Tert-amyl methyl ether | 0.51 | U | 6.0 | 0.51 |
| 637-92-3 | Ethyl t-butyl ether | 0.49 | U | 6.0 | 0.49 |
| 127-18-4 | Tetrachloroethene | 0.95 | U | 6.0 | 0.95 |
| 108-88-3 | Toluene | 0.32 | U | 6.0 | 0.32 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.38 | U | 6.0 | 0.38 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.35 | U | 6.0 | 0.35 |
| 79-01-6 | Trichloroethene | 0.45 | U | 6.0 | 0.45 |
| 75-69-4 | Trichlorofluoromethane | 0.35 | U | 6.0 | 0.35 |
| 108-05-4 | Vinyl acetate | 0.79 | U | 12 | 0.79 |
| 75-01-4 | Vinyl chloride | 0.36 | U | 6.0 | 0.36 |
| 1330-20-7 | Xylenes, Total | 0.64 | U | 12 | 0.64 |
| 98-06-6 | tert-Butylbenzene | 0.80 | U | 6.0 | 0.80 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: 146348.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/10/2015 19:01

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.63 | U | 6.0 | 0.63 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 90 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 87 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 81 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 91 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 146349.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.37(g) | Date Analyzed: 11/10/2015 19:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.55 | U | 5.5 | 0.55 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.72 | U | 5.5 | 0.72 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.34 | U | 5.5 | 0.34 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.46 | U | 5.5 | 0.46 |
| 75-34-3 | 1,1-Dichloroethane | 0.38 | U | 5.5 | 0.38 |
| 75-35-4 | 1,1-Dichloroethene | 0.89 | U | 5.5 | 0.89 |
| 563-58-6 | 1,1-Dichloropropene | 0.41 | U | 5.5 | 0.41 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.55 | U | 5.5 | 0.55 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.86 | U | 5.5 | 0.86 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.51 | U | 5.5 | 0.51 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.42 | U | 5.5 | 0.42 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.64 | U | 5.5 | 0.64 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 3.0 | U | 11 | 3.0 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.38 | U | 5.5 | 0.38 |
| 107-06-2 | 1,2-Dichloroethane | 0.51 | U | 5.5 | 0.51 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.5 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.76 | U | 5.5 | 0.76 |
| 142-28-9 | 1,3-Dichloropropane | 0.44 | U | 5.5 | 0.44 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.31 | U | 5.5 | 0.31 |
| 594-20-7 | 2,2-Dichloropropane | 0.49 | U | 5.5 | 0.49 |
| 78-93-3 | 2-Butanone (MEK) | 3.7 | J B | 22 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.71 | U | 55 | 0.71 |
| 95-49-8 | 2-Chlorotoluene | 0.33 | U | 5.5 | 0.33 |
| 591-78-6 | 2-Hexanone | 0.75 | U | 22 | 0.75 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 5.5 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.45 | U | 5.5 | 0.45 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.4 | U | 22 | 1.4 |
| 67-64-1 | Acetone | 6.3 | U | 22 | 6.3 |
| 71-43-2 | Benzene | 1.0 | U | 5.5 | 1.0 |
| 108-86-1 | Bromobenzene | 0.37 | U | 5.5 | 0.37 |
| 75-25-2 | Bromoform | 0.27 | U | 5.5 | 0.27 |
| 74-83-9 | Bromomethane | 0.46 | U | 5.5 | 0.46 |
| 75-15-0 | Carbon disulfide | 0.65 | U | 5.5 | 0.65 |
| 56-23-5 | Carbon tetrachloride | 0.72 | U | 5.5 | 0.72 |
| 108-90-7 | Chlorobenzene | 0.59 | U | 5.5 | 0.59 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 146349.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.37(g) | Date Analyzed: 11/10/2015 19:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.41 | U | 5.5 | 0.41 |
| 75-00-3 | Chloroethane | 0.49 | U | 5.5 | 0.49 |
| 67-66-3 | Chloroform | 0.40 | U | 5.5 | 0.40 |
| 74-87-3 | Chloromethane | 0.87 | U | 5.5 | 0.87 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.31 | U | 5.5 | 0.31 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.79 | U | 5.5 | 0.79 |
| 74-95-3 | Dibromomethane | 0.46 | U | 5.5 | 0.46 |
| 74-97-5 | Bromochloromethane | 0.37 | U | 5.5 | 0.37 |
| 75-71-8 | Dichlorodifluoromethane | 0.31 | U | 5.5 | 0.31 |
| 75-27-4 | Bromodichloromethane | 0.32 | U | 5.5 | 0.32 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.3 | U | 220 | 6.3 |
| 100-41-4 | Ethylbenzene | 0.35 | U | 5.5 | 0.35 |
| 87-68-3 | Hexachlorobutadiene | 0.44 | U | 5.5 | 0.44 |
| 108-20-3 | Diisopropyl ether | 0.65 | U | 11 | 0.65 |
| 98-82-8 | Isopropylbenzene | 0.22 | U | 5.5 | 0.22 |
| 1634-04-4 | Methyl tert-butyl ether | 0.48 | U | 5.5 | 0.48 |
| 75-09-2 | Methylene Chloride | 0.82 | U | 5.5 | 0.82 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.46 | U | 11 | 0.46 |
| 91-20-3 | Naphthalene | 0.39 | U | 5.5 | 0.39 |
| 104-51-8 | n-Butylbenzene | 0.65 | U | 5.5 | 0.65 |
| 103-65-1 | N-Propylbenzene | 0.66 | U | 5.5 | 0.66 |
| 95-47-6 | o-Xylene | 0.63 | U | 5.5 | 0.63 |
| 135-98-8 | sec-Butylbenzene | 1.0 | U | 5.5 | 1.0 |
| 100-42-5 | Styrene | 0.44 | U | 5.5 | 0.44 |
| 994-05-8 | Tert-amyl methyl ether | 0.48 | U | 5.5 | 0.48 |
| 637-92-3 | Ethyl t-butyl ether | 0.45 | U | 5.5 | 0.45 |
| 127-18-4 | Tetrachloroethene | 0.89 | U | 5.5 | 0.89 |
| 108-88-3 | Toluene | 0.30 | J | 5.5 | 0.30 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.35 | U | 5.5 | 0.35 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.32 | U | 5.5 | 0.32 |
| 79-01-6 | Trichloroethene | 0.42 | U | 5.5 | 0.42 |
| 75-69-4 | Trichlorofluoromethane | 0.32 | U | 5.5 | 0.32 |
| 108-05-4 | Vinyl acetate | 0.73 | U | 11 | 0.73 |
| 75-01-4 | Vinyl chloride | 0.33 | U | 5.5 | 0.33 |
| 1330-20-7 | Xylenes, Total | 0.60 | U | 11 | 0.60 |
| 98-06-6 | tert-Butylbenzene | 0.74 | U | 5.5 | 0.74 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: 146349.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.37(g) | Date Analyzed: 11/10/2015 19:22

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.59 | U | 5.5 | 0.59 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 99 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 87 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 103 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 146350.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 5.67(g) | Date Analyzed: 11/10/2015 19:44

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.52 | U | 5.2 | 0.52 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.67 | U | 5.2 | 0.67 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.32 | U | 5.2 | 0.32 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.43 | U | 5.2 | 0.43 |
| 75-34-3 | 1,1-Dichloroethane | 0.35 | U | 5.2 | 0.35 |
| 75-35-4 | 1,1-Dichloroethene | 0.83 | U | 5.2 | 0.83 |
| 563-58-6 | 1,1-Dichloropropene | 0.38 | U | 5.2 | 0.38 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.52 | U | 5.2 | 0.52 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.81 | U | 5.2 | 0.81 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.2 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.39 | U | 5.2 | 0.39 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.60 | U | 5.2 | 0.60 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.8 | U | 10 | 2.8 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.35 | U | 5.2 | 0.35 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.2 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.2 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.71 | U | 5.2 | 0.71 |
| 142-28-9 | 1,3-Dichloropropane | 0.41 | U | 5.2 | 0.41 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.29 | U | 5.2 | 0.29 |
| 594-20-7 | 2,2-Dichloropropane | 0.45 | U | 5.2 | 0.45 |
| 78-93-3 | 2-Butanone (MEK) | 3.1 | J B | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.66 | U | 52 | 0.66 |
| 95-49-8 | 2-Chlorotoluene | 0.31 | U | 5.2 | 0.31 |
| 591-78-6 | 2-Hexanone | 0.70 | U | 21 | 0.70 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.2 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.42 | U | 5.2 | 0.42 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 5.9 | U | 21 | 5.9 |
| 71-43-2 | Benzene | 0.93 | U | 5.2 | 0.93 |
| 108-86-1 | Bromobenzene | 0.34 | U | 5.2 | 0.34 |
| 75-25-2 | Bromoform | 0.25 | U | 5.2 | 0.25 |
| 74-83-9 | Bromomethane | 0.43 | U | 5.2 | 0.43 |
| 75-15-0 | Carbon disulfide | 0.61 | U | 5.2 | 0.61 |
| 56-23-5 | Carbon tetrachloride | 0.67 | U | 5.2 | 0.67 |
| 108-90-7 | Chlorobenzene | 0.55 | U | 5.2 | 0.55 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 146350.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 5.67(g) | Date Analyzed: 11/10/2015 19:44

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.38 | U | 5.2 | 0.38 |
| 75-00-3 | Chloroethane | 0.45 | U | 5.2 | 0.45 |
| 67-66-3 | Chloroform | 0.37 | U | 5.2 | 0.37 |
| 74-87-3 | Chloromethane | 0.82 | U | 5.2 | 0.82 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.29 | U | 5.2 | 0.29 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.73 | U | 5.2 | 0.73 |
| 74-95-3 | Dibromomethane | 0.43 | U | 5.2 | 0.43 |
| 74-97-5 | Bromochloromethane | 0.34 | U | 5.2 | 0.34 |
| 75-71-8 | Dichlorodifluoromethane | 0.29 | U | 5.2 | 0.29 |
| 75-27-4 | Bromodichloromethane | 0.30 | U | 5.2 | 0.30 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.9 | U | 210 | 5.9 |
| 100-41-4 | Ethylbenzene | 0.33 | U | 5.2 | 0.33 |
| 87-68-3 | Hexachlorobutadiene | 0.41 | U | 5.2 | 0.41 |
| 108-20-3 | Diisopropyl ether | 0.61 | U | 10 | 0.61 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.2 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.44 | U | 5.2 | 0.44 |
| 75-09-2 | Methylene Chloride | 0.76 | U | 5.2 | 0.76 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.43 | U | 10 | 0.43 |
| 91-20-3 | Naphthalene | 0.36 | U | 5.2 | 0.36 |
| 104-51-8 | n-Butylbenzene | 0.61 | U | 5.2 | 0.61 |
| 103-65-1 | N-Propylbenzene | 0.62 | U | 5.2 | 0.62 |
| 95-47-6 | o-Xylene | 0.59 | U | 5.2 | 0.59 |
| 135-98-8 | sec-Butylbenzene | 0.94 | U | 5.2 | 0.94 |
| 100-42-5 | Styrene | 0.41 | U | 5.2 | 0.41 |
| 994-05-8 | Tert-amyl methyl ether | 0.44 | U | 5.2 | 0.44 |
| 637-92-3 | Ethyl t-butyl ether | 0.42 | U | 5.2 | 0.42 |
| 127-18-4 | Tetrachloroethene | 0.83 | U | 5.2 | 0.83 |
| 108-88-3 | Toluene | 0.28 | U | 5.2 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.33 | U | 5.2 | 0.33 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.30 | U | 5.2 | 0.30 |
| 79-01-6 | Trichloroethene | 0.39 | U | 5.2 | 0.39 |
| 75-69-4 | Trichlorofluoromethane | 0.30 | U | 5.2 | 0.30 |
| 108-05-4 | Vinyl acetate | 0.68 | U | 10 | 0.68 |
| 75-01-4 | Vinyl chloride | 0.31 | U | 5.2 | 0.31 |
| 1330-20-7 | Xylenes, Total | 0.56 | U | 10 | 0.56 |
| 98-06-6 | tert-Butylbenzene | 0.69 | U | 5.2 | 0.69 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: 146350.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 5.67(g) | Date Analyzed: 11/10/2015 19:44

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.55 | U | 5.2 | 0.55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 105 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 98 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 90 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 100 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 146351.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.55(g) | Date Analyzed: 11/10/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.53 | U | 5.3 | 0.53 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.68 | U | 5.3 | 0.68 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.33 | U | 5.3 | 0.33 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.44 | U | 5.3 | 0.44 |
| 75-34-3 | 1,1-Dichloroethane | 0.36 | U | 5.3 | 0.36 |
| 75-35-4 | 1,1-Dichloroethene | 0.84 | U | 5.3 | 0.84 |
| 563-58-6 | 1,1-Dichloropropene | 0.39 | U | 5.3 | 0.39 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.53 | U | 5.3 | 0.53 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.82 | U | 5.3 | 0.82 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.3 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.40 | U | 5.3 | 0.40 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.61 | U | 5.3 | 0.61 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.9 | U | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.36 | U | 5.3 | 0.36 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.3 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.3 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.73 | U | 5.3 | 0.73 |
| 142-28-9 | 1,3-Dichloropropane | 0.42 | U | 5.3 | 0.42 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.30 | U | 5.3 | 0.30 |
| 594-20-7 | 2,2-Dichloropropane | 0.46 | U | 5.3 | 0.46 |
| 78-93-3 | 2-Butanone (MEK) | 4.0 | J B | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.67 | U | 53 | 0.67 |
| 95-49-8 | 2-Chlorotoluene | 0.32 | U | 5.3 | 0.32 |
| 591-78-6 | 2-Hexanone | 0.72 | U | 21 | 0.72 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.3 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 5.3 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 6.0 | U | 21 | 6.0 |
| 71-43-2 | Benzene | 0.95 | U | 5.3 | 0.95 |
| 108-86-1 | Bromobenzene | 0.35 | U | 5.3 | 0.35 |
| 75-25-2 | Bromoform | 0.25 | U | 5.3 | 0.25 |
| 74-83-9 | Bromomethane | 0.44 | U | 5.3 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.62 | U | 5.3 | 0.62 |
| 56-23-5 | Carbon tetrachloride | 0.68 | U | 5.3 | 0.68 |
| 108-90-7 | Chlorobenzene | 0.56 | U | 5.3 | 0.56 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 146351.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.55(g) | Date Analyzed: 11/10/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.39 | U | 5.3 | 0.39 |
| 75-00-3 | Chloroethane | 0.46 | U | 5.3 | 0.46 |
| 67-66-3 | Chloroform | 0.38 | U | 5.3 | 0.38 |
| 74-87-3 | Chloromethane | 0.83 | U | 5.3 | 0.83 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.30 | U | 5.3 | 0.30 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.75 | U | 5.3 | 0.75 |
| 74-95-3 | Dibromomethane | 0.44 | U | 5.3 | 0.44 |
| 74-97-5 | Bromochloromethane | 0.35 | U | 5.3 | 0.35 |
| 75-71-8 | Dichlorodifluoromethane | 0.30 | U | 5.3 | 0.30 |
| 75-27-4 | Bromodichloromethane | 0.31 | U | 5.3 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.0 | U | 210 | 6.0 |
| 100-41-4 | Ethylbenzene | 0.34 | U | 5.3 | 0.34 |
| 87-68-3 | Hexachlorobutadiene | 0.42 | U | 5.3 | 0.42 |
| 108-20-3 | Diisopropyl ether | 0.62 | U | 11 | 0.62 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.3 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.45 | U | 5.3 | 0.45 |
| 75-09-2 | Methylene Chloride | 0.78 | U | 5.3 | 0.78 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.44 | U | 11 | 0.44 |
| 91-20-3 | Naphthalene | 0.37 | U | 5.3 | 0.37 |
| 104-51-8 | n-Butylbenzene | 0.62 | U | 5.3 | 0.62 |
| 103-65-1 | N-Propylbenzene | 0.63 | U | 5.3 | 0.63 |
| 95-47-6 | o-Xylene | 0.60 | U | 5.3 | 0.60 |
| 135-98-8 | sec-Butylbenzene | 0.96 | U | 5.3 | 0.96 |
| 100-42-5 | Styrene | 0.42 | U | 5.3 | 0.42 |
| 994-05-8 | Tert-amyl methyl ether | 0.45 | U | 5.3 | 0.45 |
| 637-92-3 | Ethyl t-butyl ether | 0.43 | U | 5.3 | 0.43 |
| 127-18-4 | Tetrachloroethene | 0.84 | U | 5.3 | 0.84 |
| 108-88-3 | Toluene | 0.29 | J | 5.3 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.34 | U | 5.3 | 0.34 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.31 | U | 5.3 | 0.31 |
| 79-01-6 | Trichloroethene | 0.40 | U | 5.3 | 0.40 |
| 75-69-4 | Trichlorofluoromethane | 0.31 | U | 5.3 | 0.31 |
| 108-05-4 | Vinyl acetate | 0.70 | U | 11 | 0.70 |
| 75-01-4 | Vinyl chloride | 0.32 | U | 5.3 | 0.32 |
| 1330-20-7 | Xylenes, Total | 0.57 | U | 11 | 0.57 |
| 98-06-6 | tert-Butylbenzene | 0.71 | U | 5.3 | 0.71 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: 146351.D

Analysis Method: 8260C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.55(g) | Date Analyzed: 11/10/2015 20:05

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.56 | U | 5.3 | 0.56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 87 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 89 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 78 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 96 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 146365.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.55(g) | Date Analyzed: 11/11/2015 13:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.53 | U | 5.3 | 0.53 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.68 | U | 5.3 | 0.68 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.33 | U | 5.3 | 0.33 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.44 | U | 5.3 | 0.44 |
| 75-34-3 | 1,1-Dichloroethane | 0.36 | U | 5.3 | 0.36 |
| 75-35-4 | 1,1-Dichloroethene | 0.84 | U | 5.3 | 0.84 |
| 563-58-6 | 1,1-Dichloropropene | 0.39 | U | 5.3 | 0.39 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.53 | U | 5.3 | 0.53 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.82 | U | 5.3 | 0.82 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.48 | U | 5.3 | 0.48 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.40 | U | 5.3 | 0.40 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.61 | U | 5.3 | 0.61 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.9 | U | 11 | 2.9 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.36 | U | 5.3 | 0.36 |
| 107-06-2 | 1,2-Dichloroethane | 0.48 | U | 5.3 | 0.48 |
| 78-87-5 | 1,2-Dichloropropane | 0.15 | U | 5.3 | 0.15 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.73 | U | 5.3 | 0.73 |
| 142-28-9 | 1,3-Dichloropropane | 0.42 | U | 5.3 | 0.42 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.29 | U | 5.3 | 0.29 |
| 594-20-7 | 2,2-Dichloropropane | 0.46 | U | 5.3 | 0.46 |
| 78-93-3 | 2-Butanone (MEK) | 3.6 | J B | 21 | 1.2 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.67 | U | 53 | 0.67 |
| 95-49-8 | 2-Chlorotoluene | 0.32 | U | 5.3 | 0.32 |
| 591-78-6 | 2-Hexanone | 0.72 | U | 21 | 0.72 |
| 106-43-4 | 4-Chlorotoluene | 0.27 | U | 5.3 | 0.27 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 5.3 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.3 | U | 21 | 1.3 |
| 67-64-1 | Acetone | 6.0 | U | 21 | 6.0 |
| 71-43-2 | Benzene | 0.95 | U | 5.3 | 0.95 |
| 108-86-1 | Bromobenzene | 0.35 | U | 5.3 | 0.35 |
| 75-25-2 | Bromoform | 0.25 | U | 5.3 | 0.25 |
| 74-83-9 | Bromomethane | 0.44 | U | 5.3 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.62 | U | 5.3 | 0.62 |
| 56-23-5 | Carbon tetrachloride | 0.68 | U | 5.3 | 0.68 |
| 108-90-7 | Chlorobenzene | 0.56 | U | 5.3 | 0.56 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 146365.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.55(g) | Date Analyzed: 11/11/2015 13:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.39 | U | 5.3 | 0.39 |
| 75-00-3 | Chloroethane | 0.46 | U | 5.3 | 0.46 |
| 67-66-3 | Chloroform | 0.38 | U | 5.3 | 0.38 |
| 74-87-3 | Chloromethane | 0.83 | U | 5.3 | 0.83 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.29 | U | 5.3 | 0.29 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.75 | U | 5.3 | 0.75 |
| 74-95-3 | Dibromomethane | 0.44 | U | 5.3 | 0.44 |
| 74-97-5 | Bromochloromethane | 0.35 | U | 5.3 | 0.35 |
| 75-71-8 | Dichlorodifluoromethane | 0.29 | U | 5.3 | 0.29 |
| 75-27-4 | Bromodichloromethane | 0.31 | U | 5.3 | 0.31 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 6.0 | U | 210 | 6.0 |
| 100-41-4 | Ethylbenzene | 0.34 | U | 5.3 | 0.34 |
| 87-68-3 | Hexachlorobutadiene | 0.42 | U | 5.3 | 0.42 |
| 108-20-3 | Diisopropyl ether | 0.62 | U | 11 | 0.62 |
| 98-82-8 | Isopropylbenzene | 0.21 | U | 5.3 | 0.21 |
| 1634-04-4 | Methyl tert-butyl ether | 0.45 | U | 5.3 | 0.45 |
| 75-09-2 | Methylene Chloride | 0.78 | U | 5.3 | 0.78 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.44 | U | 11 | 0.44 |
| 91-20-3 | Naphthalene | 0.37 | U | 5.3 | 0.37 |
| 104-51-8 | n-Butylbenzene | 0.62 | U | 5.3 | 0.62 |
| 103-65-1 | N-Propylbenzene | 0.63 | U | 5.3 | 0.63 |
| 95-47-6 | o-Xylene | 0.60 | U | 5.3 | 0.60 |
| 135-98-8 | sec-Butylbenzene | 0.96 | U | 5.3 | 0.96 |
| 100-42-5 | Styrene | 0.42 | U | 5.3 | 0.42 |
| 994-05-8 | Tert-amyl methyl ether | 0.45 | U | 5.3 | 0.45 |
| 637-92-3 | Ethyl t-butyl ether | 0.43 | U | 5.3 | 0.43 |
| 127-18-4 | Tetrachloroethene | 0.84 | U | 5.3 | 0.84 |
| 108-88-3 | Toluene | 0.28 | U | 5.3 | 0.28 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.34 | U | 5.3 | 0.34 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.31 | U | 5.3 | 0.31 |
| 79-01-6 | Trichloroethene | 0.40 | U | 5.3 | 0.40 |
| 75-69-4 | Trichlorofluoromethane | 0.31 | U | 5.3 | 0.31 |
| 108-05-4 | Vinyl acetate | 0.70 | U | 11 | 0.70 |
| 75-01-4 | Vinyl chloride | 0.32 | U | 5.3 | 0.32 |
| 1330-20-7 | Xylenes, Total | 0.57 | U | 11 | 0.57 |
| 98-06-6 | tert-Butylbenzene | 0.71 | U | 5.3 | 0.71 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: 146365.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.55(g) | Date Analyzed: 11/11/2015 13:09

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.56 | U | 5.3 | 0.56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 102 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 89 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 58-123 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: TB-110515 | Lab Sample ID: 240-57597-16

Matrix: Water | Lab File ID: UXC4397.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.28 | U | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.44 | U | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.22 | U | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.30 | U | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 0.45 | U | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.37 | U | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.44 | U | 1.0 | 0.44 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.47 | U | 5.0 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.32 | U | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.41 | U | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.82 | U | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.25 | U | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 0.23 | U | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 0.25 | U | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.19 | U | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 0.19 | U | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.27 | U | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 0.26 | U | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 0.53 | U | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.59 | U | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 0.40 | U | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 0.48 | U | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 10 | 0.99 |
| 67-64-1 | Acetone | 0.94 | U | 10 | 0.94 |
| 71-43-2 | Benzene | 0.35 | U | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 0.35 | U | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 0.56 | U | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 0.44 | U | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.38 | U | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 0.43 | U | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 0.25 | U | 1.0 | 0.25 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: TB-110515 | Lab Sample ID: 240-57597-16

Matrix: Water | Lab File ID: UXC4397.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 1.0 | 0.43 |
| 75-00-3 | Chloroethane | 0.32 | U | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 0.25 | U | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 0.44 | U | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.26 | U | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.46 | U | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 0.42 | U | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.50 | U | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 4.9 | U | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 0.25 | U | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 0.35 | U | 1.0 | 0.35 |
| 108-20-3 | Diisopropyl ether | 0.50 | U | 10 | 0.50 |
| 98-82-8 | Isopropylbenzene | 0.35 | U | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 0.20 | U | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 0.33 | U | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.24 | U | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 0.45 | U | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 0.31 | U | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 0.40 | U | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 0.25 | U | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 0.48 | U | 1.0 | 0.48 |
| 100-42-5 | Styrene | 0.45 | U | 1.0 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.30 | U | 5.0 | 0.30 |
| 637-92-3 | Ethyl t-butyl ether | 0.23 | U | 5.0 | 0.23 |
| 127-18-4 | Tetrachloroethene | 0.31 | U | 1.0 | 0.31 |
| 108-88-3 | Toluene | 0.23 | U | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.30 | U | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.56 | U | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 0.22 | U | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 0.49 | U | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 0.41 | U | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 0.29 | U | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 0.52 | U | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 0.41 | U | 1.0 | 0.41 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: TB-110515 | Lab Sample ID: 240-57597-16

Matrix: Water | Lab File ID: UXC4397.D

Analysis Method: 8260C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:33

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.32 | U | 1.0 | 0.32 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 85 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 97 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 90 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 99 | | 78-125 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: P4000012.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.54(g) | Date Analyzed: 11/11/2015 13:28

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 40 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.6 | U | 40 | 9.6 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 78 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 87 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: P4000013.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:02

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.07(g) | Date Analyzed: 11/11/2015 13:45

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 39 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.4 | U | 39 | 9.4 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 39 | 16 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 84 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 100 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: P4000014.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:04

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.14(g) | Date Analyzed: 11/11/2015 14:01

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 17.8 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 15 | U | 40 | 15 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.7 | U | 40 | 9.7 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 68 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 95 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: P4000015.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:06

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.54(g) | Date Analyzed: 11/11/2015 14:17

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 40 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.6 | U | 40 | 9.6 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 75 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 104 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: P4000016.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:08

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.48(g) | Date Analyzed: 11/11/2015 14:33

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 73 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 91 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: P4000017.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:10

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.87(g) | Date Analyzed: 11/11/2015 14:50

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 40 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 40 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 40 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 40 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.6 | U | 40 | 9.6 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 40 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 40 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 76 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 109 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: P4000018.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:12

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.04(g) | Date Analyzed: 11/11/2015 15:06

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 16.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 76 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 98 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: P4000019.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:14

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.74(g) | Date Analyzed: 11/11/2015 15:23

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 18.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 15 | U | 41 | 15 |
| 11104-28-2 | Aroclor-1221 | 20 | U | 41 | 20 |
| 11141-16-5 | Aroclor-1232 | 25 | U | 41 | 25 |
| 53469-21-9 | Aroclor-1242 | 14 | U | 41 | 14 |
| 12672-29-6 | Aroclor-1248 | 9.8 | U | 41 | 9.8 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 41 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 41 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 72 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 90 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: P4000020.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:16

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.07(g) | Date Analyzed: 11/11/2015 15:39

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 19.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 15 | U | 41 | 15 |
| 11104-28-2 | Aroclor-1221 | 20 | U | 41 | 20 |
| 11141-16-5 | Aroclor-1232 | 25 | U | 41 | 25 |
| 53469-21-9 | Aroclor-1242 | 14 | U | 41 | 14 |
| 12672-29-6 | Aroclor-1248 | 9.9 | U | 41 | 9.9 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 41 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 41 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 77 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 95 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: P4000021.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:18

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 29.97(g) | Date Analyzed: 11/11/2015 15:55

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 13.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 83 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 89 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: P4000022.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:20

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.28(g) | Date Analyzed: 11/11/2015 16:12

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.5 | U | 39 | 9.5 |
| 11097-69-1 | Aroclor-1254 | 17 | U | 39 | 17 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 82 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 95 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: P4000023.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:22

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.29(g) | Date Analyzed: 11/11/2015 16:28

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 15.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 39 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 39 | 19 |
| 11141-16-5 | Aroclor-1232 | 24 | U | 39 | 24 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 39 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.4 | U | 39 | 9.4 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 39 | 16 |
| 11096-82-5 | Aroclor-1260 | 11 | U | 39 | 11 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 81 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 92 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: P4000024.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:24

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.41(g) | Date Analyzed: 11/11/2015 16:44

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 14.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 77 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 94 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: P4000025.D

Analysis Method: 8082A | Date Collected: 11/05/2015 09:26

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.19(g) | Date Analyzed: 11/11/2015 17:01

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 19 | U | 38 | 19 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.3 | U | 38 | 9.3 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 78 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 89 | | 10-155 |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: P4000029.D

Analysis Method: 8082A | Date Collected: 11/05/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30.34(g) | Date Analyzed: 11/11/2015 18:06

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 14 | U | 38 | 14 |
| 11104-28-2 | Aroclor-1221 | 18 | U | 38 | 18 |
| 11141-16-5 | Aroclor-1232 | 23 | U | 38 | 23 |
| 53469-21-9 | Aroclor-1242 | 13 | U | 38 | 13 |
| 12672-29-6 | Aroclor-1248 | 9.2 | U | 38 | 9.2 |
| 11097-69-1 | Aroclor-1254 | 16 | U | 38 | 16 |
| 11096-82-5 | Aroclor-1260 | 10 | U | 38 | 10 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 74 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 89 | | 10-155 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: A0111109.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:00

Sample wt/vol: 4.98(g) | Date Analyzed: 11/11/2015 11:08

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.7 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: A0111110.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:02

Sample wt/vol: 4.98(g) | Date Analyzed: 11/11/2015 11:45

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.0 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 79 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: A0111111.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:04

Sample wt/vol: 4.99(g) | Date Analyzed: 11/11/2015 12:23

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 17.8 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 56 | U | 120 | 56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 77 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: A0111112.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:06

Sample wt/vol: 4.99(g) | Date Analyzed: 11/11/2015 13:00

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.6 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 71 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: A0111207.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:08

Sample wt/vol: 4.98(g) | Date Analyzed: 11/12/2015 09:17

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.2 | Level: (low/med) Low

Analysis Batch No.: 206404 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 77 | J | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 71 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: A0111114.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:10

Sample wt/vol: 5.01(g) | Date Analyzed: 11/11/2015 14:13

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 76 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14  Lab Sample ID: 240-57597-7

Matrix: Solid  Lab File ID: A0111115.D

Analysis Method: 8015C  Date Collected: 11/05/2015 09:12

Sample wt/vol: 4.97(g)  Date Analyzed: 11/11/2015 14:50

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX  ID: 0.45(mm)

% Moisture: 16.2  Level: (low/med) Low

Analysis Batch No.: 206211  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: A0111116.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:14

Sample wt/vol: 4.98(g) | Date Analyzed: 11/11/2015 15:27

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 18.0 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 56 | U | 120 | 56 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 73 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: A0111119.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:16

Sample wt/vol: 5.00(g) | Date Analyzed: 11/11/2015 17:19

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 19.1 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 57 | U | 120 | 57 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: A0111120.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:18

Sample wt/vol: 5.02(g) | Date Analyzed: 11/11/2015 17:57

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 13.4 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 53 | U | 120 | 53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 79 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: A0111121.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:20

Sample wt/vol: 5.03(g) | Date Analyzed: 11/11/2015 18:34

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 16.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 55 | U | 120 | 55 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 79 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: A0111122.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:22

Sample wt/vol: 5.04(g) | Date Analyzed: 11/11/2015 19:12

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 15.9 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 78 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: A0111123.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:24

Sample wt/vol: 4.99(g) | Date Analyzed: 11/11/2015 19:48

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 14.7 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 75 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: A0111124.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:26

Sample wt/vol: 5.02(g) | Date Analyzed: 11/11/2015 20:25

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 80 | | 40-139 |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: A0111125.D

Analysis Method: 8015C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.00(g) | Date Analyzed: 11/11/2015 21:03

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 54 | U F2 | 120 | 54 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 74 | | 40-139 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-1-14 | Lab Sample ID: 240-57597-1

Matrix: Solid | Lab File ID: P5B11014.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.49(g) | Date Analyzed: 11/10/2015 20:07

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 16 | J B | 19 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 62 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-2-14 | Lab Sample ID: 240-57597-2

Matrix: Solid | Lab File ID: P5B11015.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:02

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.69(g) | Date Analyzed: 11/10/2015 20:38

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 10 | J B | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 62 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-3-14 | Lab Sample ID: 240-57597-3

Matrix: Solid | Lab File ID: P5B11016.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:04

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.32(g) | Date Analyzed: 11/10/2015 21:10

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 17.8 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.7 | U | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 59 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-4-14 | Lab Sample ID: 240-57597-4

Matrix: Solid | Lab File ID: P5B11017.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:06

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.31(g) | Date Analyzed: 11/10/2015 21:41

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.6 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.6 | J B | 20 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 60 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-5-14 | Lab Sample ID: 240-57597-5

Matrix: Solid | Lab File ID: P5B11019.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:08

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.01(g) | Date Analyzed: 11/10/2015 22:43

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.5 | U | 20 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 56 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-6-14 | Lab Sample ID: 240-57597-6

Matrix: Solid | Lab File ID: P5B11020.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:10

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.79(g) | Date Analyzed: 11/10/2015 23:14

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.7 | U | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 60 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-7-14 | Lab Sample ID: 240-57597-7

Matrix: Solid | Lab File ID: P5B11021.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:12

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.99(g) | Date Analyzed: 11/10/2015 23:46

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.2 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 16 | J B | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 65 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-8-14 | Lab Sample ID: 240-57597-8

Matrix: Solid | Lab File ID: P5B11022.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:14

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.45(g) | Date Analyzed: 11/11/2015 00:17

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 18.0 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.7 | U | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 65 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-9-14 | Lab Sample ID: 240-57597-9

Matrix: Solid | Lab File ID: P5B11023.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:16

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.38(g) | Date Analyzed: 11/11/2015 00:48

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 19.1 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.8 | U | 20 | 7.8 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 57 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-10-14 | Lab Sample ID: 240-57597-10

Matrix: Solid | Lab File ID: P5B11024.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:18

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.04(g) | Date Analyzed: 11/11/2015 01:19

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 13.4 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 13 | J B | 19 | 7.4 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 64 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-11-14 | Lab Sample ID: 240-57597-11

Matrix: Solid | Lab File ID: P5B11025.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:20

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.60(g) | Date Analyzed: 11/11/2015 01:50

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 16.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 9.0 | J B | 20 | 7.8 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 64 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-12-14 | Lab Sample ID: 240-57597-12

Matrix: Solid | Lab File ID: P5B11109.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:22

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.52(g) | Date Analyzed: 11/11/2015 19:03

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 15.9 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 30 | B | 20 | 7.7 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 73 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-13-14 | Lab Sample ID: 240-57597-13

Matrix: Solid | Lab File ID: P5B11110.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:24

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.73(g) | Date Analyzed: 11/11/2015 19:34

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 14.7 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 13 | J B | 20 | 7.6 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 68 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-14-14 | Lab Sample ID: 240-57597-14

Matrix: Solid | Lab File ID: P5B11111.D

Analysis Method: 8015C | Date Collected: 11/05/2015 09:26

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30.43(g) | Date Analyzed: 11/11/2015 20:06

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 11 | J B | 19 | 7.4 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 61 | | 40-160 |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP | Lab Sample ID: 240-57597-15

Matrix: Solid | Lab File ID: P5B11112.D

Analysis Method: 8015C | Date Collected: 11/05/2015 00:00

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 29.98(g) | Date Analyzed: 11/11/2015 20:37

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: 14.5 | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206356 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 8.2 | J B | 20 | 7.5 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 65 | | 40-160 |

**Appendix C**

Support Documentation

**LMC - BLOCK F UST**
**SDG 240-57597-1**

**F-VB-UST-2-14/F-VB-UST-DUP**

| ANALYTE | ORIGINAL | DUPLICATE | RL | RPD | RPD > 50% | ORIGINAL SAMPLE CONC >2xRL | DUPLICATE SAMPLE CONC >2xRL | DIFFERENCE >2XRL |
|---|---|---|---|---|---|---|---|---|
| BIS(2-ETHYLHEXYL)PHTHALATE* | 22 | 22 | 82 | 0.0 | FALSE | FALSE | FALSE | FALSE |

* Duplicate result is non-detected and reported at the MDL

LMC - BLOCK F UST
SDG 240-57597-1

| | |
|---|---|
| SAMPLE IDENTIFICATION | F-VB-UST-6-14 |
| COMPOUND | NAPHTHALENE |
| COMPOUND AREA | 7290 |
| INTERNAL STANDARD AMOUNT (ng) | 250 |
| DILUTION FACTOR | 1 |
| INTERNAL STANDARD AREA | 383746 |
| AVERAGE RRF | 2.3749 |
| % SOLIDS | 0.835 |
| WEIGHT OF SAMPLE (g) | 5.18 |
| ng to µg | 1000 |
| g to Kg | 1000 |
| CONCENTRATION = | 0.46 ng/g OR µg/Kg |

7290 x 250ng x 1µg x 1000g x 1/(383746 x 2.3749 x 5.18g x 0.835 x 1000ng)

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A3UX14\20151110-49103.b\146343.D
Lims ID: 240-57597-B-6-A Lab Sample ID: 240-57597-6
Client ID: F-VB-UST-6-14
Sample Type: Client
Inject. Date: 10-Nov-2015 17:13:30 ALS Bottle#: 20 Worklist Smp#: 21
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Sample Info: 240-57597-b-6-a
Misc. Info.: 240-0049103-021
Operator ID: 002808 Instrument ID: A3UX14

Method: \\ChromNA\Canton\ChromData\A3UX14\20151110-49103.b\8260_14.m
Limit Group: MSV 8260DOD ICAL
Last Update: 11-Nov-2015 03:54:33 Calib Date: 02-Sep-2015 17:24:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\ChromNA\Canton\ChromData\A3UX14\20150902-46929.b\144943.D

Column 1 : DB-624 ( 0.18 mm) Det: MS SCAN
Process Host: XAWRK007

First Level Reviewer: latat Date: 11-Nov-2015 02:56:14

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ug/l | Flags |
|---|---|---|---|---|---|---|---|---|
| * 1 Fluorobenzene | 96 | 6.641 | 6.641 | 0.000 | 100 | 1361610 | 50.0 | |
| * 2 Chlorobenzene-d5 | 117 | 9.362 | 9.362 | 0.000 | 97 | 837660 | 50.0 | |
| * 3 1,4-Dichlorobenzene-d4 | 152 | 11.338 | 11.338 | 0.000 | 98 | 414338 | 50.0 | |
| $ 4 Dibromofluoromethane (Surr | 113 | 5.966 | 5.966 | 0.000 | 94 | 262246 | 44.7 | |
| $ 5 1,2-Dichloroethane-d4 (Sur | 65 | 6.309 | 6.309 | 0.000 | 95 | 316976 | 46.3 | |
| $ 6 Toluene-d8 (Surr) | 98 | 8.120 | 8.120 | 0.000 | 94 | 1130088 | 43.8 | |
| $ 7 4-Bromofluorobenzene (Surr | 95 | 10.368 | 10.368 | 0.000 | 92 | 383746 | 40.4 | |
| 9 Dichlorodifluoromethane | 85 | | 1.316 | | | | ND | |
| 10 Chloromethane | 50 | | 1.481 | | | | ND | |
| 11 Vinyl chloride | 62 | | 1.588 | | | | ND | |
| 13 Bromomethane | 94 | | 1.919 | | | | ND | |
| 15 Chloroethane | 64 | | 2.026 | | | | ND | |
| 17 Trichlorofluoromethane | 101 | | 2.298 | | | | ND | |
| 21 1,1-Dichloroethene | 96 | | 2.878 | | | | ND | |
| 22 1,1,2-Trichloro-1,2,2-trif | 151 | | 2.913 | | | | ND | |
| 24 Acetone | 43 | | 3.008 | | | | ND | |
| 26 Carbon disulfide | 76 | | 3.126 | | | | ND | |
| 30 Methylene Chloride | 84 | | 3.576 | | | | ND | |
| 31 2-Methyl-2-propanol | 59 | | 3.860 | | | | ND | |
| 32 trans-1,2-Dichloroethene | 96 | | 3.978 | | | | ND | |
| 33 Methyl tert-butyl ether | 73 | | 4.002 | | | | ND | |
| 36 1,1-Dichloroethane | 63 | | 4.629 | | | | ND | |
| 38 Isopropyl ether | 87 | | 4.747 | | | | ND | |
| 39 Vinyl acetate | 86 | | 4.759 | | | | ND | |
| 40 Tert-butyl ethyl ether | 59 | | 5.221 | | | | ND | |
| 41 2,2-Dichloropropane | 77 | | 5.374 | | | | ND | |
| 42 cis-1,2-Dichloroethene | 96 | | 5.386 | | | | ND | |
| 43 2-Butanone (MEK) | 43 | 5.469 | 5.457 | 0.012 | 99 | 8629 | 3.29 | |
| 46 Chlorobromomethane | 128 | | 5.670 | | | | ND | |
| 49 Chloroform | 83 | | 5.789 | | | | ND | |
| 50 1,1,1-Trichloroethane | 97 | | 5.942 | | | | ND | |

| Compound | Sig | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ug/l | Flags |
|---|---|---|---|---|---|---|---|---|
| 52 Carbon tetrachloride | 117 | | 6.108 | | | | ND | |
| 53 1,1-Dichloropropene | 75 | | 6.120 | | | | ND | |
| 54 Benzene | 78 | | 6.345 | | | | ND | |
| 56 1,2-Dichloroethane | 62 | | 6.392 | | | | ND | |
| 57 Tert-amyl methyl ether | 73 | | 6.475 | | | | ND | |
| 59 Trichloroethene | 130 | | 6.996 | | | | ND | |
| 63 1,2-Dichloropropane | 63 | | 7.220 | | | | ND | |
| 65 Dibromomethane | 93 | | 7.339 | | | | ND | |
| 68 Dichlorobromomethane | 83 | | 7.493 | | | | ND | |
| 70 2-Chloroethyl vinyl ether | 63 | | 7.777 | | | | ND | |
| 71 cis-1,3-Dichloropropene | 75 | | 7.895 | | | | ND | |
| 73 4-Methyl-2-pentanone (MIBK | 43 | | 8.049 | | | | ND | |
| 74 Toluene | 91 | 8.179 | 8.179 | 0.000 | 97 | 6399 | 0.2363 | |
| 75 trans-1,3-Dichloropropene | 75 | | 8.392 | | | | ND | |
| 78 Tetrachloroethene | 164 | | 8.640 | | | | ND | |
| 79 1,3-Dichloropropane | 76 | | 8.699 | | | | ND | |
| 80 2-Hexanone | 43 | | 8.770 | | | | ND | |
| 82 Chlorodibromomethane | 129 | | 8.889 | | | | ND | |
| 83 Ethylene Dibromide | 107 | | 8.983 | | | | ND | |
| 85 Chlorobenzene | 112 | | 9.398 | | | | ND | |
| 86 1,1,1,2-Tetrachloroethane | 131 | | 9.469 | | | | ND | |
| 87 Ethylbenzene | 106 | | 9.480 | | | | ND | |
| 88 m-Xylene & p-Xylene | 106 | | 9.575 | | | | ND | |
| 89 o-Xylene | 106 | | 9.918 | | | | ND | |
| 90 Styrene | 104 | | 9.930 | | | | ND | |
| 91 Bromoform | 173 | | 10.108 | | | | ND | |
| 92 Isopropylbenzene | 105 | | 10.226 | | | | ND | |
| 94 Bromobenzene | 156 | | 10.498 | | | | ND | |
| 95 1,1,2,2-Tetrachloroethane | 83 | | 10.498 | | | | ND | |
| 96 1,2,3-Trichloropropane | 110 | | 10.534 | | | | ND | |
| 98 N-Propylbenzene | 120 | | 10.569 | | | | ND | |
| 99 2-Chlorotoluene | 126 | | 10.652 | | | | ND | |
| 101 4-Chlorotoluene | 126 | | 10.747 | | | | ND | |
| 103 tert-Butylbenzene | 119 | | 10.995 | | | | ND | |
| 105 1,2,4-Trimethylbenzene | 105 | | 11.031 | | | | ND | |
| 107 sec-Butylbenzene | 105 | | 11.173 | | | | ND | |
| 108 1,3-Dichlorobenzene | 146 | | 11.279 | | | | ND | |
| 109 4-Isopropyltoluene | 119 | | 11.303 | | | | ND | |
| 110 1,4-Dichlorobenzene | 146 | | 11.362 | | | | ND | |
| 111 1,2,3-Trimethylbenzene | 105 | | 11.397 | | | | ND | |
| 113 n-Butylbenzene | 91 | | 11.646 | | | | ND | |
| 114 1,2-Dichlorobenzene | 146 | | 11.670 | | | | ND | |
| 115 1,2-Dibromo-3-Chloropropan | 157 | | 12.344 | | | | ND | |
| 117 1,2,4-Trichlorobenzene | 180 | 13.018 | 13.019 | 0.000 | 33 | 1666 | 0.2066 | |
| 118 Hexachlorobutadiene | 225 | | 13.137 | | | | ND | |
| 119 Naphthalene | 128 | 13.220 | 13.220 | 0.000 | 95 | 7290 | 0.3888 | |
| 120 1,2,3-Trichlorobenzene | 180 | 13.409 | 13.409 | 0.000 | 83 | 1768 | 0.2313 | |
| S 134 Xylenes, Total | 106 | | 16.530 | | | | ND | |

LMC - BLOCK F UST AREA
SDG 240-57597-1

| | |
|---|---|
| SAMPLE IDENTIFICATION | F-VB-UST-13-14 |
| COMPOUND | BIS(2-ETHYLHEXYL)PHTHALATE |
| COMPOUND AREA | 108069 |
| INTERNAL STANDARD AMOUNT (ng) | 4 |
| DILUTION FACTOR | 1 |
| INTERNAL STANDARD AREA | 785246 |
| AVERAGE RRF | 0.7086 |
| VOLUME FINAL EXTRACT (µl) | 2000 |
| WEIGHT OF SAMPLE (g) | 30.37 |
| VOLUME INJECTED (µl) | 1 |
| ng to g | 1000 |
| g to Kg | 1000 |
| PERCENT SOLIDS | 0.8530 |

CONCENTRATION = 60 µg/Kg

108069 x 4ng x 2000µl x 1000g x 1µg x 1/(785246 x 0.7086 x 30.37g x 0.853 x 1µl x 1Kg x 1000ng)

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\A4HP9\20151111-49136.b\51111021.D
Lims ID: 240-57597-E-13-B Lab Sample ID: 240-57597-13
Client ID: F-VB-UST-13-14
Sample Type: Client
Inject. Date: 11-Nov-2015 15:39:30 ALS Bottle#: 21 Worklist Smp#: 21
Injection Vol: 1.0 ul Dil. Factor: 1.0000
Sample Info: 240-0049136-021
Misc. Info.: 240-57597-E-13-B
Operator ID: 001710 Instrument ID: A4HP9

Method: \\ChromNA\Canton\ChromData\A4HP9\20151111-49136.b\8270 _9.m
Limit Group: MSS 8270D ICAL
Last Update: 12-Nov-2015 09:01:43 Calib Date: 30-Oct-2015 16:42:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\ChromNA\Canton\ChromData\A4HP9\20151030-48759.b\51030010.D

Column 1 : 5% phenyl ( 0.18 mm) Det: MS SCAN
Process Host: XAWRK012

First Level Reviewer: ulmanm Date: 12-Nov-2015 08:13:03

| Compound | Sig | RT (min.) | Adj RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng/ul | Flags |
|---|---|---|---|---|---|---|---|---|
| * 1 1,4-Dichlorobenzene-d4 | 152 | 6.295 | 6.298 | -0.003 | 97 | 229993 | 4.00 | |
| * 2 Naphthalene-d8 | 136 | 7.416 | 7.418 | -0.002 | 100 | 1005170 | 4.00 | |
| * 3 Acenaphthene-d10 | 164 | 8.921 | 8.924 | -0.003 | 92 | 509477 | 4.00 | |
| * 4 Phenanthrene-d10 | 188 | 10.186 | 10.188 | -0.002 | 97 | 754672 | 4.00 | |
| * 5 Chrysene-d12 | 240 | 12.455 | 12.458 | -0.003 | 96 | 785246 | 4.00 | |
| * 6 Perylene-d12 | 264 | 14.177 | 14.184 | -0.007 | 96 | 684137 | 4.00 | |
| $ 7 2-Fluorophenol | 112 | 5.059 | 5.043 | 0.016 | 96 | 565296 | 6.51 | |
| $ 8 Phenol-d5 | 99 | 5.959 | 5.947 | 0.012 | 98 | 644249 | 5.71 | |
| $ 9 Nitrobenzene-d5 | 82 | 6.781 | 6.779 | 0.002 | 87 | 761421 | 6.46 | |
| $ 10 2-Fluorobiphenyl (Surr) | 172 | 8.320 | 8.322 | -0.002 | 99 | 944233 | 5.88 | |
| $ 11 2,4,6-Tribromophenol | 330 | 9.589 | 9.592 | -0.003 | 81 | 59786 | 3.35 | M |
| $ 12 Terphenyl-d14 | 244 | 11.503 | 11.506 | -0.003 | 99 | 1112006 | 7.75 | |
| 41 Benzaldehyde | 77 | | 5.894 | | | | ND | |
| 43 Phenol | 94 | | 5.961 | | | | ND | |
| 69 Bis(2-chloroethyl)ether | 93 | | 6.033 | | | | ND | |
| 46 2-Chlorophenol | 128 | | 6.115 | | | | ND | |
| 51 2-Methylphenol | 108 | | 6.500 | | | | ND | |
| 53 2,2'-oxybis[1-chloropropan | 45 | | 6.519 | | | | ND | |
| 57 N-Nitrosodi-n-propylamine | 70 | | 6.630 | | | | ND | |
| 21 3 & 4 Methylphenol | 108 | | 6.630 | | | | ND | |
| 58 Acetophenone | 105 | | 6.644 | | | | ND | |
| 61 Hexachloroethane | 117 | | 6.760 | | | | ND | |
| 62 Nitrobenzene | 77 | | 6.798 | | | | ND | |
| 64 Isophorone | 82 | | 7.000 | | | | ND | |
| 66 2-Nitrophenol | 139 | | 7.077 | | | | ND | |
| 65 2,4-Dimethylphenol | 107 | | 7.091 | | | | ND | |
| 70 Bis(2-chloroethoxy)methane | 93 | | 7.163 | | | | ND | |
| 73 2,4-Dichlorophenol | 162 | | 7.289 | | | | ND | |
| 75 Naphthalene | 128 | | 7.438 | | | | ND | |
| 76 4-Chloroaniline | 127 | | 7.462 | | | | ND | |
| 79 Hexachlorobutadiene | 225 | | 7.539 | | | | ND | |

| Compound | Sig | RT (min.) | Adj RT (min.) | Dlt RT (min.) | Q | Response | OnCol Amt ng/ul | Flags |
|---|---|---|---|---|---|---|---|---|
| 84 Caprolactam | 113 | | 7.750 | | | | ND | |
| 86 4-Chloro-3-methylphenol | 107 | | 7.861 | | | | ND | |
| 87 2-Methylnaphthalene | 142 | | 8.024 | | | | ND | |
| 89 Hexachlorocyclopentadiene | 237 | | 8.164 | | | | ND | |
| 92 2,4,6-Trichlorophenol | 196 | | 8.260 | | | | ND | |
| 93 2,4,5-Trichlorophenol | 196 | | 8.298 | | | | ND | |
| 98 1,1'-Biphenyl | 154 | | 8.414 | | | | ND | |
| 100 2-Chloronaphthalene | 162 | | 8.447 | | | | ND | |
| 102 2-Nitroaniline | 65 | | 8.510 | | | | ND | |
| 105 Dimethyl phthalate | 163 | | 8.640 | | | | ND | |
| 107 2,6-Dinitrotoluene | 165 | | 8.702 | | | | ND | |
| 109 Acenaphthylene | 152 | | 8.803 | | | | ND | |
| 110 3-Nitroaniline | 138 | | 8.856 | | | | ND | |
| 111 2,4-Dinitrophenol | 184 | | 8.943 | | | | ND | |
| 112 Acenaphthene | 153 | | 8.948 | | | | ND | |
| 113 4-Nitrophenol | 109 | | 8.981 | | | | ND | |
| 114 2,4-Dinitrotoluene | 165 | | 9.049 | | | | ND | |
| 116 Dibenzofuran | 168 | | 9.092 | | | | ND | |
| 121 Diethyl phthalate | 149 | | 9.226 | | | | ND | |
| 123 4-Chlorophenyl phenyl ethe | 204 | | 9.356 | | | | ND | |
| 125 4-Nitroaniline | 138 | | 9.376 | | | | ND | |
| 126 Fluorene | 166 | | 9.385 | | | | ND | |
| 128 4,6-Dinitro-2-methylphenol | 198 | | 9.404 | | | | ND | |
| 131 N-Nitrosodiphenylamine | 169 | | 9.452 | | | | ND | |
| 139 4-Bromophenyl phenyl ether | 248 | | 9.775 | | | | ND | |
| 141 Hexachlorobenzene | 284 | | 9.866 | | | | ND | |
| 142 Atrazine | 200 | | 9.876 | | | | ND | |
| 143 Pentachlorophenol | 266 | | 10.020 | | | | ND | |
| 148 Phenanthrene | 178 | | 10.207 | | | | ND | |
| 149 Anthracene | 178 | | 10.251 | | | | ND | |
| 151 Carbazole | 167 | | 10.371 | | | | ND | |
| 154 Di-n-butyl phthalate | 149 | 10.599 | 10.602 | -0.003 | 99 | 51250 | 0.2300 | |
| 159 Fluoranthene | 202 | | 11.222 | | | | ND | |
| 161 Pyrene | 202 | | 11.424 | | | | ND | |
| 168 Butyl benzyl phthalate | 149 | | 11.876 | | | | ND | |
| 173 Bis(2-ethylhexyl) phthalat | 149 | 12.345 | 12.347 | -0.002 | 95 | 108069 | 0.7769 | |
| 175 3,3'-Dichlorobenzidine | 252 | | 12.386 | | | | ND | |
| 177 Benzo[a]anthracene | 228 | | 12.448 | | | | ND | |
| 178 Chrysene | 228 | | 12.487 | | | | ND | |
| 180 Di-n-octyl phthalate | 149 | | 12.982 | | | | ND | |
| 182 Benzo[b]fluoranthene | 252 | | 13.646 | | | | ND | |
| 183 Benzo[k]fluoranthene | 252 | | 13.679 | | | | ND | |
| 184 Benzo[a]pyrene | 252 | | 14.103 | | | | ND | |
| 188 Indeno[1,2,3-cd]pyrene | 276 | | 16.060 | | | | ND | |
| 189 Dibenz(a,h)anthracene | 278 | | 16.065 | | | | ND | |
| 190 Benzo[g,h,i]perylene | 276 | | 16.646 | | | | ND | |

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\AFID\20151112-49174.b\A0111207.D
Lims ID: 240-57597-C-5-B Lab Sample ID: 240-57597-5
Client ID: F-VB-UST-5-14
Sample Type: Client
Inject. Date: 12-Nov-2015 09:17:06 ALS Bottle#: 7 Worklist Smp#: 7
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Sample Info:
Operator ID: Instrument ID: AFID

Method: \\ChromNA\Canton\ChromData\AFID\20151112-49174.b\A80218015.m
Limit Group: GCV 8015C GRO ICAL
Last Update: 12-Nov-2015 09:30:49 Calib Date: 24-Aug-2015 17:07:27
Integrator: Falcon
Quant Method: External Standard Quant By: Initial Calibration
Last ICal File: \\ChromNA\Canton\ChromData\AFID\20150824-46596.b\A0082112.D

Column 1 : Det: GC FID1A
Process Host: XAWRK031

| Compound | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Response | OnCol Amt ug/l | Flags |
|---|---|---|---|---|---|---|
| $ 16 Trifluorotoluene (Surr) | 10.372 | 10.368 | 0.004 | 3169615 | 7.13 | |
| A 5 C6-C10 | 11.532 | (5.929-17.135) | | 26342362 | 65.1 | |

## Reagents:

VGGASURROGATE_00034 Amount Added: 1.00 Units: uL Run Reagent

4.98g/5ml
% solids = 84.8%
CF = 404923.648
DF = 1

26342362/404923.648 ug/L x 0.005L/4.98g x 1/0.848 x 1000g/1Kg
= 77 ug/Kg

0.8/C0.9
2.1/C2.2

# Chain of Custody Record #46 016125

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING
TestAmerica Laboratories, Inc.
TAL-8210 (0713)

Regulatory Program: ☐ DW ☐ NPDES ☐ RCRA ☐ Other:

| Client Contact | Project Manager: SCOTT NESBIT | Site Contact: J. MULLIS | Date: 11/5/15 | COC No: |
|---|---|---|---|---|
| Company Name: TETRA TECH | Tel/Fax: 412-559-8133 | Lab Contact: J. MCFADDEN | Carrier: FedEx | 1 of 2 COCs |
| Address: 20251 CENTURY BLVD STE 200 | Analysis Turnaround Time | | | Sampler: |
| City/State/Zip: GERMANTOWN, MD, 20874 | ☐ CALENDAR DAYS ☐ WORKING DAYS | | | For Lab Use Only: |
| Phone: 301-528-3000 | TAT if different from Below _____ | | | Walk-in Client: |
| Fax: 301-528-3000 | ☐ 2 weeks | | | Lab Sampling: |
| Project Name: MRC SOIL REMENT | ☒ 1 week | | | |
| Site: BLOCK F UST | ☐ 2 days | | | Job / SDG No.: |
| PO # 112IC07336 | ☐ 1 day | | | |

| Sample Identification | Sample Date | Sample Time | Sample Type (C=Comp, G=Grab) | Matrix | # of Cont. | Filtered Sample (Y/N) | Perform MS / MSD (Y/N) | VOCs (8260C) | SVOCs (8270D) | TPH GRO/DRO (8015B) | TCLP Metals (1311 then 6010C) | PCBs (8082A) | Ignitability | Sample Specific Notes: |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| F-VB-UST-1-14 | 11/5/15 | 0900 | G | SO | 4 | | | X | X | X | X | X | | |
| F-VB-UST-2-14 | | 0902 | | | | | | X | X | X | X | X | | |
| F-VB-UST-3-14 | | 0904 | | | | | | X | X | X | X | X | | |
| F-VB-UST-4-14 | | 0906 | | | | | | X | X | X | X | X | | |
| F-VB-UST-5-14 | | 0908 | | | | | | X | X | X | X | X | | |
| F-VB-UST-6-14 | | 0910 | | | | | | X | X | X | X | X | | |
| F-VB-UST-7-14 | | 0912 | | | | | | X | X | X | X | X | | |
| F-VB-UST-8-14 | | 0914 | | | | | | X | X | X | X | X | | |
| F-VB-UST-9-14 | | 0916 | | | | | | X | X | X | X | X | | |
| F-VB-UST-10-14 | | 0918 | | | | | | X | X | X | X | X | | |
| F-VB-UST-11-14 | | 0920 | | | | | | X | X | X | X | X | | |
| F-VB-UST-12-14 | | 0922 | | | | | | X | X | X | X | X | | |

Preservation Used: 1= Ice, 2= HCl, 3= H2SO4, 4=HNO3, 5=NaOH, 6= Other _____

**Possible Hazard Identification:**
Are any samples from a listed EPA Hazardous Waste? Please List any EPA Waste Codes for the sample in the Comments Section if the lab is to dispose of the sample.

☒ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

**Sample Disposal ( A fee may be assessed if samples are retained longer than 1 month)**

☐ Return to Client ☒ Disposal by Lab ☐ Archive for _____ Months

**Special Instructions/QC Requirements & Comments:** MRC SOIL REMENT: BLOCK F UST BASE CONFIRMATION SAMPLES

| Custody Seals Intact: ☐ Yes ☐ No | Custody Seal No.: | | Cooler Temp. (°C): Obs'd: _____ Corr'd: _____ | | Therm ID No.: _____ |
|---|---|---|---|---|---|
| Relinquished by: | Company: TETRA TECH | Date/Time: 11/5/15 1530 | Received by: | Company: TestAmerica | Date/Time: 11/5/15 1530 |
| Relinquished by: | Company: TestAmerica | Date/Time: 11/5/15 1618 | Received by: | Company: TA | Date/Time: 11.6.15 1020 |
| Relinquished by: | Company: | Date/Time: | Received in Laboratory by: | Company: | Date/Time: |

# Chain of Custody Record

#46

016124

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING
TestAmerica Laboratories, Inc.
TAL-8210 (0713)

Regulatory Program: ☐ DW ☐ NPDES ☐ RCRA ☐ Other:

| Client Contact | Project Manager: SCOTT NESBIT | Site Contact: J. Mullis | Date: 11/5/15 | COC No: |
|---|---|---|---|---|
| Company Name: TETRA TECH | Tel/Fax: 412-559-8133 | Lab Contact: J. McFadden | Carrier: FedEx | 2 of 2 COCs |
| Address: 20251 Century Blvd. Ste. 200 | Analysis Turnaround Time | | | Sampler: |
| City/State/Zip: Germantown, MD, 20874 | ☐ CALENDAR DAYS ☐ WORKING DAYS | | | For Lab Use Only: |
| Phone: 301-528-3000 | TAT if different from Below _____ | | | Walk-in Client: |
| Fax: 301-528-3000 | ☐ 2 weeks | | | Lab Sampling: |
| Project Name: MRC SOIL REMEDY | ⊗ 1 week | | | |
| Site: BLOCK F UST CONFIRMATION | ☐ 2 days | | | Job / SDG No.: |
| P O # 112E07336 | ☐ 1 day | | | |

| Sample Identification | Sample Date | Sample Time | Sample Type (C=Comp, G=Grab) | Matrix | # of Cont. | Filtered Sample (Y/N) | Perform MS / MSD (Y/N) | VOCs (8260C) | SVOCs (8270D) | TPH Gro/Dro (8015B) | TCLP Metals (1311+6010) | RCRA (8015B) | Sample Specific Notes: |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| F-VB-UST-13-14 | 11/5/15 | 0924 | G | SO | 4 | | | X | X | X | X | X | |
| F-VB-UST-14-14 | | 0926 | | | | | | X | X | X | X | X | |
| F-VB-UST-DUP | | — | | | | | | X | X | X | X | X | # DUPLICATE |
| TB-110515 | | — | | AQ | 2 | | | X | | | | | |

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11/5/15

Preservation Used: (1)= Ice, 2= HCl, 3= H2SO4, 4=HNO3, 5=NaOH, 6= Other

**Possible Hazard Identification:**
Are any samples from a listed EPA Hazardous Waste? Please List any EPA Waste Codes for the sample in the Comments Section if the lab is to dispose of the sample.

☒ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

**Sample Disposal ( A fee may be assessed if samples are retained longer than 1 month)**

☐ Return to Client ☒ Disposal by Lab ☐ Archive for _____ Months

**Special Instructions/QC Requirements & Comments:** SEE PAGE 1

| Custody Seals Intact: ☐ Yes ☐ No | Custody Seal No.: | | Cooler Temp. (°C): Obs'd: _____ Corr'd: _____ Therm ID No.: _____ | | |
|---|---|---|---|---|---|
| Relinquished by: | Company: TETRA TECH | Date/Time: 11/5/15 1530 | Received by: | Company: TestAmerica | Date/Time: 11/5/15 1530 |
| Relinquished by: | Company: TestAmerica | Date/Time: 11/5/15 1618 | Received by: | Company: TA | Date/Time: 11-6-15 1020 |
| Relinquished by: | Company: | Date/Time: | Received in Laboratory by: | Company: | Date/Time: |

**TestAmerica Canton Sample Receipt Form/Narrative** Login # : 57507
**Canton Facility**

Client Tetra Tech Site Name _____ Cooler unpacked by:

Cooler Received on 11-6-15 Opened on 11-6-15

FedEx: 1st Grd ~~Exp~~ UPS FAS Stetson Client Drop Off TestAmerica Courier Other _____

**Receipt After-hours**: Drop-off Date/Time _____ Storage Location _____

TestAmerica Cooler # _____ Foam Box Client Cooler Box Other _____

Packing material used: ~~Bubble Wrap~~ Foam Plastic Bag None Other _____

COOLANT: ~~Wet Ice~~ Blue Ice Dry Ice Water None

1. Cooler temperature upon receipt

IR GUN# 53 (CF **+0.1** °C) Observed Cooler Temp._____°C Corrected Cooler Temp._____°C

IR GUN# 48 (CF **-0.3** °C) Observed Cooler Temp._____°C Corrected Cooler Temp._____°C ☑ See Multiple Cooler Form

~~IR GUN# 5 (CF +0.4 °C) Observed Cooler Temp._____°C Corrected Cooler Temp._____°C~~

IR GUN# 8 (CF **-0.5** °C) Observed Cooler Temp._____°C Corrected Cooler Temp._____°C

| | | |
|---|---|---|
| 2. Were custody seals on the outside of the cooler(s)? If Yes Quantity 2 | ~~Yes~~ | No |
| -Were custody seals on the outside of the cooler(s) signed & dated? | ~~Yes~~ | No NA |
| -Were custody seals on the bottle(s) or bottle kits (LLHg/MeHg)? | Yes | ~~No~~ |
| 3. Shippers' packing slip attached to the cooler(s)? | ~~Yes~~ | No |
| 4. Did custody papers accompany the sample(s)? | ~~Yes~~ | No |
| 5. Were the custody papers relinquished & signed in the appropriate place? | ~~Yes~~ | No |
| 6. Was/were the person(s) who collected the samples clearly identified on the COC? | Yes | ~~No~~ |
| 7. Did all bottles arrive in good condition (Unbroken)? | ~~Yes~~ | No |
| 8. Could all bottle labels be reconciled with the COC? | ~~Yes~~ | No |
| 9. Were correct bottle(s) used for the test(s) indicated? | ~~Yes~~ | No |
| 10. Sufficient quantity received to perform indicated analyses? | ~~Yes~~ | No |
| 11. Were sample(s) at the correct pH upon receipt? | Yes No | ~~NA~~ pH Strip Lot# **HC554612** |
| 12. Were VOAs on the COC? | ~~Yes~~ | No |
| 13. Were air bubbles >6 mm in any VOA vials? | Yes | ~~No~~ NA |
| 14. Was a trip blank present in the cooler(s)? Trip Blank Lot #_____ | ~~Yes~~ | No |

Contacted PM _____ Date _____ by _____ via Verbal Voice Mail Other

Concerning _____

**14. CHAIN OF CUSTODY & SAMPLE DISCREPANCIES** Samples processed by:

**15. SAMPLE CONDITION**

Sample(s) _____ were received after the recommended holding time had expired.

Sample(s) _____ were received in a broken container.

Sample(s) _____ were received with bubble >6 mm in diameter. (Notify PM)

**16. SAMPLE PRESERVATION**

Sample(s) _____ were further preserved in the laboratory.

Time preserved: _____ Preservative(s) added/Lot number(s): _____

TestAmerica Multiple Cooler Receipt Form/Narrative
Canton Facility

Login # : 575597

| Cooler # | IR Gun # | Observed Temp °C | Corrected Temp °C | Coolant |
|---|---|---|---|---|
| Client | 53 | 0.8 | 0.9 | ice |
| 2 | 1 | 2.1 | 2.2 | 1 |
| | | | | |
| | | | | |
| | | | |

# CASE NARRATIVE

## Client: Tetra Tech, Inc.

## Project: MRC Soil Remedy Block F UST

## Report Number: 240-57597-1

With the exceptions noted as flags or footnotes, standard analytical protocols were followed in the analysis of the samples and no problems were encountered or anomalies observed. In addition all laboratory quality control samples were within established control limits, with any exceptions noted below. Each sample was analyzed to achieve the lowest possible reporting limit within the constraints of the method. In some cases, due to interference or analytes present at high concentrations, samples were diluted. For diluted samples, the reporting limits are adjusted relative to the dilution required.

TestAmerica Canton attests to the validity of the laboratory data generated by TestAmerica facilities reported herein. All analyses performed by TestAmerica facilities were done using established laboratory SOPs that incorporate QA/QC procedures described in the application methods. TestAmerica's operations groups have reviewed the data for compliance with the laboratory QA/QC plan, and data have been found to be compliant with laboratory protocols unless otherwise noted below.

The test results in this report meet all NELAP requirements for parameters for which accreditation is required or available. Any exceptions to NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory.

Calculations are performed before rounding to avoid round-off errors in calculated results.

All holding times were met and proper preservation noted for the methods performed on these samples, unless otherwise detailed in the individual sections below.

All solid sample results are reported on an "as received" basis unless otherwise indicated by the presence of a % solids value in the method header.

This laboratory report is confidential and is intended for the sole use of TestAmerica and its client.

**RECEIPT**

The samples were received on 11/6/2015 10:20 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperatures of the 2 coolers at receipt time were 0.9º C and 2.2º C.

**VOLATILE ORGANIC COMPOUNDS (GCMS)**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for volatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8260C. The samples were prepared and analyzed on 11/10/2015 and 11/11/2015.

2-Butanone (MEK), Naphthalene and Toluene were detected in method blank MB 240-206267/6 at levels that were above the method detection limit but below the reporting limit. The values should be considered estimates, and have been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

2-Butanone (MEK) was detected in method blank MB 240-206089/1-A at a level that was above the method detection limit but below the reporting limit. The value should be considered an estimate, and has been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

The continuing calibration verification (CCV) associated with batch 240-206051 recovered above the upper control limit for 1,1,1-Trichloroethane; 1,1,2-Trichloro-1,2,2-trifluoroethane; 2,2-Dichloropropane; Carbon tetrachloride; Dichloro-difluoromethane and Trichlorofluoromethane. The samples associated with this CCV were non-detects for the affected analytes; therefore, the data have been reported.

The continuing calibration verification (CCV) analyzed in batch 240-206267 was outside the method criteria for the following analyte(s): 2-Hexanone, Acetone, 2-Methyl-2-propanol, 2-Chloroethyl Vinyl Ether, 4-Methyl-2-Pentanone, 2,2-Dichloropropane, and 2-Butanone. An MRL standard at or below the reporting limit (RL) was analyzed with the affected samplesF-VB-UST-DUP (240-57597-15), (CCV 240-206267/4), (CCVIS 240-206267/3), (LCS 240-206267/5), (MB 240-206267/6), (240-57597-B-15-C MS) and (240-57597-B-15-D MSD) and found to be acceptable. As indicated in the reference method, sample analysis may proceed; however, any detection for the affected analyte(s) is considered estimated.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**VOLATILE ORGANIC COMPOUNDS (GCMS)**

Sample TB-110515 (240-57597-16) was analyzed for volatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8260C. The samples were analyzed on 11/12/2015.

2-Butanone (MEK) was detected in method blank MB 240-206461/6 at a level that was above the method detection limit but below the reporting limit. The value should be considered an estimate, and has been flagged. If the associated sample reported a result above the MDL and/or RL, the result has been flagged.

The continuing calibration verification (CCV) for analytical batch 206461 exceeded control criteria for Hexachlorobutadiene. The samples associated with this CCV were non-detect for the affected analyte. In accordance with the laboratory SOP, a low level CCV at the reporting limit (labeled as an MRL) was analyzed and the affected compound was detected; therefore the data has been reported. No further corrective action was required. TB-110515 (240-57597-16)

2-Chloroethyl vinyl ether cannot be reliably recovered in an acid preserved sample. TB-110515 (240-57597-16)

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**SEMIVOLATILE ORGANIC COMPOUNDS (GCMS)**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for semivolatile organic compounds (GCMS) in accordance with EPA SW-846 Method 8270D. The samples were prepared on 11/09/2015 and analyzed on 11/11/2015.

The analyte benzaldehyde had an Rf below method recommendations in analytical batch 206233. An LODV was analyzed to demonstrate sensitivity for this analyte at the reporting limit for the following samples:F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14), F-VB-UST-DUP (240-57597-15), (LCS 240-205866/9-A), (MB 240-205866/8-A), (240-57597-D-15-E MS) and (240-57597-D-15-F MSD).

Two analytes had Rfs below method recommendations: Benzaldehyde and Hexachlorocyclopentadiene. A LODV was analyzed to demonstrate sensitivity for these analytes at the reporting limit for the following samples: F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9) and F-VB-UST-10-14 (240-57597-10).

The continuing calibration verification (CCV) associated with analytical batch 206233 recovered above the upper control limit for N-Nitrosodi-n-propylamine. The samples associated with this CCV were non-detect for this analyte; therefore, the data have been reported. The following samples are impacted: F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14), F-VB-UST-DUP (240-57597-15), (MB 240-205866/8-A), (240-57597-D-15-E MS) and (240-57597-D-15-F MSD).

The continuing calibration verification (CCV) for analytical batch 206233 was below method recommendations for the analyte benzaldehyde (>20%D). Benzaldehyde was present in sample F-VB-UST-11-14 (240-57597-11) at an estimated amount of 19ug/kg J. Had the response of benzaldehyde been within method recommendations the amount of benzaldehyde measured in the sample would likely have still been a small estimated concentration (J flag). Therefore, no additional analysis or corrective action was initiated. The results have been flagged accordingly and reported.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**GASOLINE RANGE ORGANICS**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for gasoline range organics in accordance with EPA SW-846 Method 8015C - GRO. The samples were prepared and analyzed on 11/11/2015 and 11/12/2015.

The matrix spike duplicate (MSD) precision for 206211 was outside control limits. Sample matrix interference and/or non-homogeneity are suspected because the associated laboratory control sample (LCS) precision was within acceptance limits.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**DIESEL RANGE ORGANICS (DRO)**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12),

F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for diesel range organics (DRO) in accordance with EPA SW-846 Method 8015C - DRO. The samples were prepared on 11/09/2015 and analyzed on 11/10/2015 and 11/11/2015.

Total Petroleum Hydrocarbons (C10-C32) was detected in method blank MB 240-205788/22-A at a level exceeding the reporting limit. If the associated sample reported a result above the MDL and/or RL, the result has been flagged. .

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**POLYCHLORINATED BIPHENYLS (PCBS)**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for polychlorinated biphenyls (PCBs) in accordance with EPA SW-846 Method 8082A. The samples were prepared on 11/09/2015 and analyzed on 11/11/2015.

Surrogates are added during the extraction process prior to dilution. When the sample dilution is 5X or greater, surrogate recoveries are diluted out and no corrective action is required. All of the samples in this data set analyzed for PCBs were subjected to the sulfuric acid cleanup procedure before instrumental analysis, per EPA Method 3665A.

The following samples required a tetrabutylammonium sulfite (TBA) clean-up to reduce matrix interferences caused by sulfur: F-VB-UST-5-14 (240-57597-5) and F-VB-UST-13-14 (240-57597-13).Reagents 2266653, 1848564 and 2283702.

No additional analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

**PERCENT SOLIDS**

Samples F-VB-UST-1-14 (240-57597-1), F-VB-UST-2-14 (240-57597-2), F-VB-UST-3-14 (240-57597-3), F-VB-UST-4-14 (240-57597-4), F-VB-UST-5-14 (240-57597-5), F-VB-UST-6-14 (240-57597-6), F-VB-UST-7-14 (240-57597-7), F-VB-UST-8-14 (240-57597-8), F-VB-UST-9-14 (240-57597-9), F-VB-UST-10-14 (240-57597-10), F-VB-UST-11-14 (240-57597-11), F-VB-UST-12-14 (240-57597-12), F-VB-UST-13-14 (240-57597-13), F-VB-UST-14-14 (240-57597-14) and F-VB-UST-DUP (240-57597-15) were analyzed for percent solids in accordance with EPA Method 160.3 MOD. The samples were analyzed on 11/09/2015.

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

## Qualifiers

### GC/MS VOA

| Qualifier | Qualifier Description |
|---|---|
| U | Indicates the analyte was analyzed for but not detected. |
| B | Compound was found in the blank and sample. |
| J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| F2 | MS/MSD RPD exceeds control limits |
| F1 | MS and/or MSD Recovery is outside acceptance limits. |

### GC/MS Semi VOA

| Qualifier | Qualifier Description |
|---|---|
| U | Indicates the analyte was analyzed for but not detected. |
| J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |

### GC VOA

| Qualifier | Qualifier Description |
|---|---|
| U | Indicates the analyte was analyzed for but not detected. |
| F2 | MS/MSD RPD exceeds control limits |
| J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |

### GC Semi VOA

| Qualifier | Qualifier Description |
|---|---|
| B | Compound was found in the blank and sample. |
| J | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |
| U | Indicates the analyte was analyzed for but not detected. |

## Glossary

| Abbreviation | These commonly used abbreviations may or may not be present in this report. |
|---|---|
| ¤ | Listed under the "D" column to designate that the result is reported on a dry weight basis |
| %R | Percent Recovery |
| CFL | Contains Free Liquid |
| CNF | Contains no Free Liquid |
| DER | Duplicate error ratio (normalized absolute difference) |
| Dil Fac | Dilution Factor |
| DL, RA, RE, IN | Indicates a Dilution, Re-analysis, Re-extraction, or additional Initial metals/anion analysis of the sample |
| DLC | Decision level concentration |
| MDA | Minimum detectable activity |
| EDL | Estimated Detection Limit |
| MDC | Minimum detectable concentration |
| MDL | Method Detection Limit |
| ML | Minimum Level (Dioxin) |
| NC | Not Calculated |
| ND | Not detected at the reporting limit (or MDL or EDL if shown) |
| PQL | Practical Quantitation Limit |
| QC | Quality Control |
| RER | Relative error ratio |
| RL | Reporting Limit or Requested Limit (Radiochemistry) |
| RPD | Relative Percent Difference, a measure of the relative difference between two points |
| TEF | Toxicity Equivalent Factor (Dioxin) |
| TEQ | Toxicity Equivalent Quotient (Dioxin) |

# Sample Summary

Client: Tetra Tech, Inc.
Project/Site: MRC Soil Remedy Block F UST

TestAmerica Job ID: 240-57597-1

| Lab Sample ID | Client Sample ID | Matrix | Collected | Received |
|---|---|---|---|---|
| 240-57597-1 | F-VB-UST-1-14 | Solid | 11/05/15 09:00 | 11/06/15 10:20 |
| 240-57597-2 | F-VB-UST-2-14 | Solid | 11/05/15 09:02 | 11/06/15 10:20 |
| 240-57597-3 | F-VB-UST-3-14 | Solid | 11/05/15 09:04 | 11/06/15 10:20 |
| 240-57597-4 | F-VB-UST-4-14 | Solid | 11/05/15 09:06 | 11/06/15 10:20 |
| 240-57597-5 | F-VB-UST-5-14 | Solid | 11/05/15 09:08 | 11/06/15 10:20 |
| 240-57597-6 | F-VB-UST-6-14 | Solid | 11/05/15 09:10 | 11/06/15 10:20 |
| 240-57597-7 | F-VB-UST-7-14 | Solid | 11/05/15 09:12 | 11/06/15 10:20 |
| 240-57597-8 | F-VB-UST-8-14 | Solid | 11/05/15 09:14 | 11/06/15 10:20 |
| 240-57597-9 | F-VB-UST-9-14 | Solid | 11/05/15 09:16 | 11/06/15 10:20 |
| 240-57597-10 | F-VB-UST-10-14 | Solid | 11/05/15 09:18 | 11/06/15 10:20 |
| 240-57597-11 | F-VB-UST-11-14 | Solid | 11/05/15 09:20 | 11/06/15 10:20 |
| 240-57597-12 | F-VB-UST-12-14 | Solid | 11/05/15 09:22 | 11/06/15 10:20 |
| 240-57597-13 | F-VB-UST-13-14 | Solid | 11/05/15 09:24 | 11/06/15 10:20 |
| 240-57597-14 | F-VB-UST-14-14 | Solid | 11/05/15 09:26 | 11/06/15 10:20 |
| 240-57597-15 | F-VB-UST-DUP | Solid | 11/05/15 00:00 | 11/06/15 10:20 |
| 240-57597-16 | TB-110515 | Water | 11/05/15 00:00 | 11/06/15 10:20 |

# Method Summary

Client: Tetra Tech, Inc.
Project/Site: MRC Soil Remedy Block F UST

TestAmerica Job ID: 240-57597-1

| Method | Method Description | Protocol | Laboratory |
|---|---|---|---|
| 8260C | Volatile Organic Compounds by GC/MS | SW846 | TAL CAN |
| 8270D | Semivolatile Organic Compounds (GC/MS) | SW846 | TAL CAN |
| 8015C | Nonhalogenated Organics using GC/FID -Modified (Gasoline Range Organics) | SW846 | TAL CAN |
| 8015C | Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics) | SW846 | TAL CAN |
| 8082A | Polychlorinated Biphenyls (PCBs) by Gas Chromatography | SW846 | TAL CAN |
| Moisture | Percent Moisture | EPA | TAL CAN |

**Protocol References:**

EPA = US Environmental Protection Agency

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

**Laboratory References:**

TAL CAN = TestAmerica Canton, 4101 Shuffel Street NW, North Canton, OH 44720, TEL (330)497-9396

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): DB-624 ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | DBFM # | DCA # | TOL # | BFB # |
|---|---|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 101 | 103 | 103 | 91 |
| F-VB-UST-2-14 | 240-57597-2 | 95 | 99 | 96 | 85 |
| F-VB-UST-3-14 | 240-57597-3 | 90 | 90 | 93 | 86 |
| F-VB-UST-4-14 | 240-57597-4 | 94 | 102 | 93 | 83 |
| F-VB-UST-5-14 | 240-57597-5 | 101 | 104 | 100 | 87 |
| F-VB-UST-6-14 | 240-57597-6 | 89 | 93 | 88 | 81 |
| F-VB-UST-7-14 | 240-57597-7 | 98 | 103 | 98 | 87 |
| F-VB-UST-8-14 | 240-57597-8 | 93 | 93 | 97 | 85 |
| F-VB-UST-9-14 | 240-57597-9 | 88 | 94 | 89 | 82 |
| F-VB-UST-10-14 | 240-57597-10 | 88 | 93 | 89 | 83 |
| F-VB-UST-11-14 | 240-57597-11 | 87 | 91 | 90 | 81 |
| F-VB-UST-12-14 | 240-57597-12 | 98 | 103 | 99 | 87 |
| F-VB-UST-13-14 | 240-57597-13 | 98 | 100 | 105 | 90 |
| F-VB-UST-14-14 | 240-57597-14 | 89 | 96 | 87 | 78 |
| F-VB-UST-DUP | 240-57597-15 | 101 | 102 | 102 | 89 |
| | MB 240-206089/1-A | 88 | 91 | 91 | 82 |
| | MB 240-206267/6 | 101 | 102 | 101 | 90 |
| | LCS 240-206051/5 | 95 | 92 | 94 | 87 |
| | LCS 240-206267/5 | 99 | 94 | 95 | 89 |
| F-VB-UST-5-14 MS | 240-57597-5 MS | 97 | 92 | 99 | 95 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 97 | 97 | 97 | 88 |
| F-VB-UST-5-14 MSD | 240-57597-5 MSD | 101 | 98 | 101 | 96 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 99 | 95 | 96 | 86 |

| | QC LIMITS |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 37-132 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 58-123 |
| TOL = Toluene-d8 (Surr) | 67-125 |
| BFB = 4-Bromofluorobenzene (Surr) | 52-136 |

# Column to be used to flag recovery values

FORM II 8260C

FORM II
GC/MS VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low

GC Column (1): DB-624 ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | DBFM # | DCA # | TOL # | BFB # |
|---|---|---|---|---|---|
| TB-110515 | 240-57597-16 | 97 | 99 | 85 | 90 |
| | MB 240-206461/6 | 99 | 95 | 88 | 97 |
| | LCS 240-206461/4 | 103 | 99 | 96 | 103 |
| | 240-57564-B-1 MS | 99 | 97 | 88 | 103 |
| | 240-57564-B-1 MSD | 101 | 98 | 90 | 102 |

| | QC LIMITS |
|---|---|
| DBFM = Dibromofluoromethane (Surr) | 79-120 |
| DCA = 1,2-Dichloroethane-d4 (Surr) | 78-125 |
| TOL = Toluene-d8 (Surr) | 80-120 |
| BFB = 4-Bromofluorobenzene (Surr) | 61-120 |

# Column to be used to flag recovery values

FORM II 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: ICIS 240-196026/6  Date Analyzed: 09/02/2015 15:37

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 144938.D  Heated Purge: (Y/N) Y

Calibration ID: 30511

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 1489699 | 6.64 | 910038 | 9.37 | 406298 | 11.34 |
| UPPER LIMIT | | 2979398 | 7.14 | 1820076 | 9.87 | 812596 | 11.84 |
| LOWER LIMIT | | 744850 | 6.14 | 455019 | 8.87 | 203149 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCVIS 240-206051/3 | | 1498068 | 6.64 | 927026 | 9.36 | 430150 | 11.34 |
| CCVIS 240-206267/3 | | 1511907 | 6.64 | 908062 | 9.37 | 416832 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVIS 240-206051/3  Date Analyzed: 11/10/2015 10:47

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 146325.D  Heated Purge: (Y/N) Y

Calibration ID: 30511

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1498068 | 6.64 | 927026 | 9.36 | 430150 | 11.34 |
| UPPER LIMIT | | 2996136 | 7.14 | 1854052 | 9.86 | 860300 | 11.84 |
| LOWER LIMIT | | 749034 | 6.14 | 463513 | 8.86 | 215075 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-206051/4 | | 1428310 | 6.64 | 843706 | 9.36 | 400898 | 11.34 |
| LCS 240-206051/5 | | 1546829 | 6.64 | 933267 | 9.37 | 439911 | 11.34 |
| MB 240-206089/1-A | | 1539926 | 6.64 | 908904 | 9.36 | 440491 | 11.34 |
| 240-57597-1 | F-VB-UST-1-14 | 1339894 | 6.64 | 803741 | 9.36 | 401204 | 11.34 |
| 240-57597-2 | F-VB-UST-2-14 | 1366253 | 6.64 | 812014 | 9.36 | 402470 | 11.34 |
| 240-57597-3 | F-VB-UST-3-14 | 1478321 | 6.64 | 869963 | 9.36 | 422508 | 11.34 |
| 240-57597-4 | F-VB-UST-4-14 | 1379125 | 6.64 | 832253 | 9.36 | 407647 | 11.34 |
| 240-57597-5 | F-VB-UST-5-14 | 1381401 | 6.64 | 834119 | 9.36 | 405215 | 11.34 |
| 240-57597-5 MS | F-VB-UST-5-14 MS | 1327604 | 6.64 | 776997 | 9.36 | 334424 | 11.34 |
| 240-57597-5 MSD | F-VB-UST-5-14 MSD | 1417706 | 6.64 | 855433 | 9.36 | 375351 | 11.34 |
| 240-57597-6 | F-VB-UST-6-14 | 1361610 | 6.64 | 837660 | 9.36 | 414338 | 11.34 |
| 240-57597-7 | F-VB-UST-7-14 | 1397574 | 6.64 | 848234 | 9.36 | 420468 | 11.34 |
| 240-57597-8 | F-VB-UST-8-14 | 1221245 | 6.64 | 712201 | 9.36 | 345698 | 11.34 |
| 240-57597-9 | F-VB-UST-9-14 | 1423543 | 6.64 | 857394 | 9.36 | 412539 | 11.34 |
| 240-57597-10 | F-VB-UST-10-14 | 1539979 | 6.64 | 947956 | 9.36 | 457583 | 11.34 |
| 240-57597-11 | F-VB-UST-11-14 | 1382482 | 6.64 | 822070 | 9.36 | 396271 | 11.34 |
| 240-57597-12 | F-VB-UST-12-14 | 1354784 | 6.64 | 808462 | 9.36 | 408687 | 11.34 |
| 240-57597-13 | F-VB-UST-13-14 | 1299068 | 6.64 | 749701 | 9.36 | 370193 | 11.34 |
| 240-57597-14 | F-VB-UST-14-14 | 1212543 | 6.64 | 728498 | 9.36 | 358599 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVIS 240-206267/3  Date Analyzed: 11/11/2015 10:39

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 146358.D  Heated Purge: (Y/N) Y

Calibration ID: 30511

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1511907 | 6.64 | 908062 | 9.37 | 416832 | 11.34 |
| UPPER LIMIT | | 3023814 | 7.14 | 1816124 | 9.87 | 833664 | 11.84 |
| LOWER LIMIT | | 755954 | 6.14 | 454031 | 8.87 | 208416 | 10.84 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-206267/4 | | 1401932 | 6.64 | 838774 | 9.36 | 393967 | 11.34 |
| LCS 240-206267/5 | | 1488652 | 6.64 | 913203 | 9.36 | 430177 | 11.34 |
| MB 240-206267/6 | | 1376673 | 6.64 | 826304 | 9.37 | 402736 | 11.34 |
| 240-57597-15 | F-VB-UST-DUP | 1258982 | 6.64 | 751722 | 9.36 | 367644 | 11.34 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 953829 | 6.64 | 561900 | 9.36 | 271317 | 11.34 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 1326891 | 6.64 | 791618 | 9.36 | 374548 | 11.34 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: STD8260 240-204067/4  Date Analyzed: 10/28/2015 13:51

Instrument ID: A3UX15  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UXC3826A.D  Heated Purge: (Y/N) N

Calibration ID: 31159

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 2262723 | 4.98 | 1708800 | 7.64 | 900225 | 9.85 |
| UPPER LIMIT | | 4525446 | 5.48 | 3417600 | 8.14 | 1800450 | 10.35 |
| LOWER LIMIT | | 1131362 | 4.48 | 854400 | 7.14 | 450113 | 9.35 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-204067/14 | | 2244648 | 4.98 | 1717477 | 7.64 | 915948 | 9.85 |
| ICV 240-204067/15 | | 2191668 | 4.97 | 1675976 | 7.64 | 861710 | 9.85 |
| CCVIS 240-206461/2 | | 1685909 | 4.97 | 1402013 | 7.64 | 853216 | 9.85 |

FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM VIII
GC/MS VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVIS 240-206461/2  Date Analyzed: 11/12/2015 10:41

Instrument ID: A3UX15  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): UXC4392.D  Heated Purge: (Y/N) N

Calibration ID: 31161

| | | FB | | CBZ | | DCB | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 1685909 | 4.97 | 1402013 | 7.64 | 853216 | 9.85 |
| UPPER LIMIT | | 3371818 | 5.47 | 2804026 | 8.14 | 1706432 | 10.35 |
| LOWER LIMIT | | 842955 | 4.47 | 701007 | 7.14 | 426608 | 9.35 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| LCS 240-206461/4 | | 1696051 | 4.97 | 1338674 | 7.64 | 737315 | 9.85 |
| CCV 240-206461/3 | | 1607750 | 4.97 | 1329730 | 7.64 | 731454 | 9.85 |
| MB 240-206461/6 | | 1530091 | 4.97 | 1276242 | 7.64 | 702637 | 9.85 |
| 240-57597-16 | TB-110515 | 1547502 | 4.97 | 1309915 | 7.64 | 691488 | 9.85 |
| 240-57564-B-1 MS | | 1604429 | 4.97 | 1330852 | 7.64 | 805667 | 9.85 |
| 240-57564-B-1 MSD | | 1611108 | 4.97 | 1343144 | 7.64 | 800978 | 9.85 |

FB = Fluorobenzene
FB = Fluorobenzene
CBZ = Chlorobenzene-d5
DCB = 1,4-Dichlorobenzene-d4
Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8260C

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab File ID: 146361.D  Lab Sample ID: MB 240-206267/6

Matrix: Solid  Heated Purge:(Y/N) Y

Instrument ID: A3UX14  Date Analyzed: 11/11/2015 11:43

GC Column: DB-624  ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-206267/5 | 146360.D | 11/11/2015 11:22 |
| F-VB-UST-DUP | 240-57597-15 | 146365.D | 11/11/2015 13:09 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 146366.D | 11/11/2015 13:31 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 146367.D | 11/11/2015 13:52 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206267/6

Matrix: Solid | Lab File ID: 146361.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/11/2015 11:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.80 | U | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.78 | U | 5.0 | 0.78 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 3.14 | J | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.90 | U | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206267/6

Matrix: Solid | Lab File ID: 146361.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/11/2015 11:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.79 | U | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.71 | U | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.7 | U | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.74 | U | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.350 | J | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.91 | U | 5.0 | 0.91 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.80 | U | 5.0 | 0.80 |
| 108-88-3 | Toluene | 0.369 | J | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206267/6

Matrix: Solid | Lab File ID: 146361.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(g) | Date Analyzed: 11/11/2015 11:43

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206267 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 101 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 101 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 90 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 102 | | 58-123 |

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146360.D

Lab ID: LCS 240-206267/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 21.7 | 108 | 80-120 | |
| 1,1,1-Trichloroethane | 20.0 | 22.4 | 112 | 77-126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 20.2 | 101 | 77-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 21.2 | 106 | 80-138 | |
| 1,1-Dichloroethane | 20.0 | 19.0 | 95 | 76-120 | |
| 1,1-Dichloroethene | 20.0 | 18.6 | 93 | 75-135 | |
| 1,1-Dichloropropene | 20.0 | 20.9 | 104 | 80-120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 20.6 | 103 | 61-121 | |
| 1,2,3-Trichloropropane | 20.0 | 21.8 | 109 | 80-126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 21.3 | 106 | 64-124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 21.1 | 106 | 80-129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 20.9 | 104 | 61-132 | |
| 1,2-Dichlorobenzene | 20.0 | 20.6 | 103 | 76-120 | |
| 1,2-Dichloroethane | 20.0 | 20.6 | 103 | 72-120 | |
| 1,2-Dichloropropane | 20.0 | 19.1 | 96 | 80-120 | |
| 1,3-Dichlorobenzene | 20.0 | 20.2 | 101 | 78-120 | |
| 1,3-Dichloropropane | 20.0 | 20.5 | 102 | 80-120 | |
| 1,4-Dichlorobenzene | 20.0 | 20.4 | 102 | 75-120 | |
| 2,2-Dichloropropane | 20.0 | 21.9 | 110 | 69-135 | |
| 2-Butanone (MEK) | 40.0 | 33.7 | 84 | 52-131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 15.3 J | 76 | 28-173 | |
| 2-Chlorotoluene | 20.0 | 21.5 | 108 | 78-120 | |
| 2-Hexanone | 40.0 | 32.9 | 82 | 64-136 | |
| 4-Chlorotoluene | 20.0 | 21.6 | 108 | 79-120 | |
| p-Isopropyltoluene | 20.0 | 22.3 | 112 | 77-131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 32.6 | 82 | 67-135 | |
| Acetone | 40.0 | 33.3 | 83 | 41-137 | |
| Benzene | 20.0 | 19.4 | 97 | 79-120 | |
| Bromobenzene | 20.0 | 22.4 | 112 | 80-120 | |
| Bromoform | 20.0 | 17.7 | 89 | 62-133 | |
| Bromomethane | 20.0 | 19.6 | 98 | 42-136 | |
| Carbon disulfide | 20.0 | 14.8 | 74 | 62-146 | |
| Carbon tetrachloride | 20.0 | 22.8 | 114 | 71-129 | |
| Chlorobenzene | 20.0 | 21.0 | 105 | 78-120 | |
| Dibromochloromethane | 20.0 | 19.4 | 97 | 72-127 | |
| Chloroethane | 20.0 | 19.9 | 99 | 58-120 | |
| Chloroform | 20.0 | 21.0 | 105 | 77-120 | |
| Chloromethane | 20.0 | 15.6 | 78 | 50-120 | |
| cis-1,2-Dichloroethene | 20.0 | 20.1 | 101 | 76-120 | |
| cis-1,3-Dichloropropene | 20.0 | 18.4 | 92 | 74-128 | |
| Dibromomethane | 20.0 | 21.4 | 107 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146360.D

Lab ID: LCS 240-206267/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 20.0 | 22.2 | 111 | 79-120 | |
| Dichlorodifluoromethane | 20.0 | 20.4 | 102 | 26-120 | |
| Bromodichloromethane | 20.0 | 19.2 | 96 | 80-122 | |
| tert-Butyl alcohol (TBA) | 200 | 143 J | 71 | 60-125 | |
| Ethylbenzene | 20.0 | 21.3 | 106 | 79-120 | |
| Hexachlorobutadiene | 20.0 | 23.6 | 118 | 54-131 | |
| Isopropylbenzene | 20.0 | 21.2 | 106 | 76-122 | |
| Methyl tert-butyl ether | 20.0 | 18.5 | 93 | 49-165 | |
| Methylene Chloride | 20.0 | 16.9 | 84 | 75-120 | |
| m-Xylene & p-Xylene | 20.0 | 21.3 | 107 | 80-120 | |
| Naphthalene | 20.0 | 18.4 | 92 | 65-123 | |
| n-Butylbenzene | 20.0 | 21.8 | 109 | 68-135 | |
| N-Propylbenzene | 20.0 | 21.7 | 108 | 80-129 | |
| o-Xylene | 20.0 | 21.7 | 108 | 80-120 | |
| sec-Butylbenzene | 20.0 | 20.9 | 105 | 74-129 | |
| Styrene | 20.0 | 20.0 | 100 | 80-120 | |
| Tetrachloroethene | 20.0 | 21.3 | 107 | 79-120 | |
| Toluene | 20.0 | 21.3 | 106 | 75-120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.3 | 101 | 78-120 | |
| trans-1,3-Dichloropropene | 20.0 | 20.6 | 103 | 73-131 | |
| Trichloroethene | 20.0 | 21.5 | 108 | 79-120 | |
| Trichlorofluoromethane | 20.0 | 22.8 | 114 | 57-146 | |
| Vinyl acetate | 20.0 | 17.7 | 88 | 70-130 | |
| Vinyl chloride | 20.0 | 16.8 | 84 | 57-120 | |
| Xylenes, Total | 40.0 | 43.0 | 108 | 80-120 | |
| tert-Butylbenzene | 20.0 | 21.4 | 107 | 76-126 | |
| 1,2-Dibromoethane | 20.0 | 21.7 | 109 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146366.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 19.8 | 0.53 U | 20.2 | 102 | 34-135 | |
| 1,1,1-Trichloroethane | 19.8 | 0.68 U | 22.0 | 111 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 19.8 | 0.33 U | 19.2 | 97 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 19.8 | 0.44 U | 21.4 | 108 | 50-147 | |
| 1,1-Dichloroethane | 19.8 | 0.36 U | 18.4 | 93 | 54-122 | |
| 1,1-Dichloroethene | 19.8 | 0.84 U | 18.3 | 93 | 49-157 | |
| 1,1-Dichloropropene | 19.8 | 0.39 U | 20.4 | 103 | 50-122 | |
| 1,2,3-Trichlorobenzene | 19.8 | 0.53 U | 19.1 | 96 | 10-120 | |
| 1,2,3-Trichloropropane | 19.8 | 0.82 U | 22.1 | 112 | 32-174 | |
| 1,2,4-Trichlorobenzene | 19.8 | 0.40 U | 19.5 | 98 | 10-120 | |
| 1,2,4-Trimethylbenzene | 19.8 | 0.61 U | 19.4 | 98 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 19.8 | 2.9 U | 17.8 | 90 | 10-153 | |
| 1,2-Dichlorobenzene | 19.8 | 0.36 U | 18.5 | 94 | 17-122 | |
| 1,2-Dichloroethane | 19.8 | 0.48 U | 19.3 | 97 | 49-123 | |
| 1,2-Dichloropropane | 19.8 | 0.15 U | 17.0 | 86 | 61-120 | |
| 1,3-Dichlorobenzene | 19.8 | 0.73 U | 19.2 | 97 | 16-126 | |
| 1,3-Dichloropropane | 19.8 | 0.42 U | 18.5 | 94 | 54-128 | |
| 1,4-Dichlorobenzene | 19.8 | 0.29 U | 17.9 | 91 | 15-121 | |
| 2,2-Dichloropropane | 19.8 | 0.46 U | 21.6 | 109 | 49-132 | |
| 2-Butanone (MEK) | 39.6 | 3.6 J | 30.1 | 67 | 30-143 | |
| 2-Chloroethyl vinyl ether | 19.8 | 0.67 U | 12.6 J | 63 | 10-162 | |
| 2-Chlorotoluene | 19.8 | 0.32 U | 19.9 | 101 | 11-162 | |
| 2-Hexanone | 39.6 | 0.72 U | 31.1 | 79 | 37-147 | |
| 4-Chlorotoluene | 19.8 | 0.27 U | 19.3 | 98 | 15-149 | |
| p-Isopropyltoluene | 19.8 | 0.43 U | 20.1 | 102 | 10-165 | |
| 4-Methyl-2-pentanone (MIBK) | 39.6 | 1.3 U | 32.3 | 82 | 43-147 | |
| Acetone | 39.6 | 6.0 U | 30.0 | 76 | 24-140 | |
| Benzene | 19.8 | 0.95 U | 18.5 | 94 | 53-120 | |
| Bromobenzene | 19.8 | 0.35 U | 19.9 | 101 | 24-144 | |
| Bromoform | 19.8 | 0.25 U | 16.1 | 81 | 18-129 | |
| Bromomethane | 19.8 | 0.44 U | 19.1 | 97 | 33-130 | |
| Carbon disulfide | 19.8 | 0.62 U | 13.8 | 70 | 20-151 | |
| Carbon tetrachloride | 19.8 | 0.68 U | 21.4 | 108 | 32-137 | |
| Chlorobenzene | 19.8 | 0.56 U | 19.3 | 97 | 37-120 | |
| Dibromochloromethane | 19.8 | 0.39 U | 17.1 | 86 | 29-135 | |
| Chloroethane | 19.8 | 0.46 U | 19.4 | 98 | 45-120 | |
| Chloroform | 19.8 | 0.38 U | 20.5 | 104 | 53-120 | |
| Chloromethane | 19.8 | 0.83 U | 15.5 | 78 | 34-120 | |
| cis-1,2-Dichloroethene | 19.8 | 0.29 U | 19.6 | 99 | 50-120 | |
| cis-1,3-Dichloropropene | 19.8 | 0.75 U | 16.0 | 81 | 27-133 | |
| Dibromomethane | 19.8 | 0.44 U | 21.4 | 108 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146366.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Bromochloromethane | 19.8 | 0.35 U | 20.6 | 104 | 53-120 | |
| Dichlorodifluoromethane | 19.8 | 0.29 U | 19.9 | 101 | 17-120 | |
| Bromodichloromethane | 19.8 | 0.31 U | 17.2 | 87 | 35-132 | |
| tert-Butyl alcohol (TBA) | 198 | 6.0 U | 163 J | 83 | 60-125 | |
| Ethylbenzene | 19.8 | 0.34 U | 19.0 | 96 | 30-131 | |
| Hexachlorobutadiene | 19.8 | 0.42 U | 20.9 | 106 | 10-131 | |
| Isopropylbenzene | 19.8 | 0.21 U | 20.0 | 101 | 21-134 | |
| Methyl tert-butyl ether | 19.8 | 0.45 U | 17.5 | 88 | 51-157 | |
| Methylene Chloride | 19.8 | 0.78 U | 16.2 | 82 | 54-120 | |
| m-Xylene & p-Xylene | 19.8 | 0.44 U | 19.4 | 98 | 29-131 | |
| Naphthalene | 19.8 | 0.37 U | 16.9 | 85 | 10-124 | |
| n-Butylbenzene | 19.8 | 0.62 U | 19.3 | 98 | 10-148 | |
| N-Propylbenzene | 19.8 | 0.63 U | 19.4 | 98 | 10-178 | |
| o-Xylene | 19.8 | 0.60 U | 19.5 | 98 | 29-134 | |
| sec-Butylbenzene | 19.8 | 0.96 U | 19.5 | 98 | 10-172 | |
| Styrene | 19.8 | 0.42 U | 17.8 | 90 | 27-127 | |
| Tetrachloroethene | 19.8 | 0.84 U | 21.7 | 109 | 31-135 | |
| Toluene | 19.8 | 0.28 U | 20.2 | 102 | 39-129 | |
| trans-1,2-Dichloroethene | 19.8 | 0.34 U | 19.7 | 100 | 50-123 | |
| trans-1,3-Dichloropropene | 19.8 | 0.31 U | 17.1 | 86 | 28-137 | |
| Trichloroethene | 19.8 | 0.40 U | 20.3 | 103 | 10-177 | |
| Trichlorofluoromethane | 19.8 | 0.31 U | 21.2 | 107 | 36-142 | |
| Vinyl acetate | 19.8 | 0.70 U | 17.6 | 89 | 70-130 | |
| Vinyl chloride | 19.8 | 0.32 U | 16.9 | 85 | 42-120 | |
| Xylenes, Total | 39.6 | 0.57 U | 38.9 | 98 | 30-131 | |
| tert-Butylbenzene | 19.8 | 0.71 U | 20.0 | 101 | 10-163 | |
| 1,2-Dibromoethane | 19.8 | 0.56 U | 19.8 | 100 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146367.D

Lab ID: 240-57597-15 MSD Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 22.6 | 22.0 | 97 | 9 | 30 | 34-135 | |
| 1,1,1-Trichloroethane | 22.6 | 23.8 | 105 | 8 | 30 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 22.6 | 21.3 | 94 | 11 | 30 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 22.6 | 23.5 | 104 | 10 | 30 | 50-147 | |
| 1,1-Dichloroethane | 22.6 | 20.4 | 90 | 10 | 30 | 54-122 | |
| 1,1-Dichloroethene | 22.6 | 20.3 | 90 | 10 | 30 | 49-157 | |
| 1,1-Dichloropropene | 22.6 | 22.2 | 98 | 8 | 30 | 50-122 | |
| 1,2,3-Trichlorobenzene | 22.6 | 19.9 | 88 | 4 | 30 | 10-120 | |
| 1,2,3-Trichloropropane | 22.6 | 20.5 | 91 | 8 | 30 | 32-174 | |
| 1,2,4-Trichlorobenzene | 22.6 | 20.8 | 92 | 6 | 30 | 10-120 | |
| 1,2,4-Trimethylbenzene | 22.6 | 20.8 | 92 | 7 | 30 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 22.6 | 20.5 | 91 | 14 | 30 | 10-153 | |
| 1,2-Dichlorobenzene | 22.6 | 19.9 | 88 | 7 | 30 | 17-122 | |
| 1,2-Dichloroethane | 22.6 | 21.1 | 93 | 9 | 30 | 49-123 | |
| 1,2-Dichloropropane | 22.6 | 19.2 | 85 | 12 | 30 | 61-120 | |
| 1,3-Dichlorobenzene | 22.6 | 19.0 | 84 | 1 | 30 | 16-126 | |
| 1,3-Dichloropropane | 22.6 | 20.2 | 89 | 9 | 30 | 54-128 | |
| 1,4-Dichlorobenzene | 22.6 | 19.1 | 85 | 7 | 30 | 15-121 | |
| 2,2-Dichloropropane | 22.6 | 24.1 | 107 | 11 | 30 | 49-132 | |
| 2-Butanone (MEK) | 45.2 | 31.6 | 62 | 5 | 30 | 30-143 | |
| 2-Chloroethyl vinyl ether | 22.6 | 14.8 J | 66 | 17 | 30 | 10-162 | |
| 2-Chlorotoluene | 22.6 | 21.8 | 96 | 9 | 30 | 11-162 | |
| 2-Hexanone | 45.2 | 33.3 | 74 | 7 | 30 | 37-147 | |
| 4-Chlorotoluene | 22.6 | 20.1 | 89 | 4 | 30 | 15-149 | |
| p-Isopropyltoluene | 22.6 | 21.4 | 94 | 6 | 30 | 10-165 | |
| 4-Methyl-2-pentanone (MIBK) | 45.2 | 33.2 | 73 | 3 | 30 | 43-147 | |
| Acetone | 45.2 | 29.1 | 64 | 3 | 30 | 24-140 | |
| Benzene | 22.6 | 20.6 | 91 | 11 | 30 | 53-120 | |
| Bromobenzene | 22.6 | 22.2 | 98 | 11 | 30 | 24-144 | |
| Bromoform | 22.6 | 17.5 | 77 | 8 | 30 | 18-129 | |
| Bromomethane | 22.6 | 20.8 | 92 | 8 | 30 | 33-130 | |
| Carbon disulfide | 22.6 | 15.4 | 68 | 11 | 30 | 20-151 | |
| Carbon tetrachloride | 22.6 | 23.7 | 105 | 10 | 30 | 32-137 | |
| Chlorobenzene | 22.6 | 20.8 | 92 | 8 | 30 | 37-120 | |
| Dibromochloromethane | 22.6 | 19.4 | 86 | 13 | 30 | 29-135 | |
| Chloroethane | 22.6 | 20.8 | 92 | 7 | 30 | 45-120 | |
| Chloroform | 22.6 | 22.6 | 100 | 9 | 30 | 53-120 | |
| Chloromethane | 22.6 | 17.2 | 76 | 11 | 30 | 34-120 | |
| cis-1,2-Dichloroethene | 22.6 | 21.5 | 95 | 9 | 30 | 50-120 | |
| cis-1,3-Dichloropropene | 22.6 | 18.1 | 80 | 12 | 30 | 27-133 | |
| Dibromomethane | 22.6 | 22.8 | 101 | 7 | 30 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 146367.D

Lab ID: 240-57597-15 MSD  Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Bromochloromethane | 22.6 | 22.2 | 98 | 8 | 30 | 53-120 | |
| Dichlorodifluoromethane | 22.6 | 21.8 | 96 | 9 | 30 | 17-120 | |
| Bromodichloromethane | 22.6 | 20.4 | 90 | 17 | 30 | 35-132 | |
| tert-Butyl alcohol (TBA) | 226 | 158 J | 70 | 4 | 30 | 60-125 | |
| Ethylbenzene | 22.6 | 20.9 | 93 | 10 | 30 | 30-131 | |
| Hexachlorobutadiene | 22.6 | 22.4 | 99 | 7 | 30 | 10-131 | |
| Isopropylbenzene | 22.6 | 21.8 | 96 | 8 | 30 | 21-134 | |
| Methyl tert-butyl ether | 22.6 | 19.5 | 86 | 11 | 30 | 51-157 | |
| Methylene Chloride | 22.6 | 19.2 | 85 | 17 | 30 | 54-120 | |
| m-Xylene & p-Xylene | 22.6 | 20.2 | 89 | 4 | 30 | 29-131 | |
| Naphthalene | 22.6 | 18.8 | 83 | 11 | 30 | 10-124 | |
| n-Butylbenzene | 22.6 | 20.7 | 92 | 7 | 30 | 10-148 | |
| N-Propylbenzene | 22.6 | 20.7 | 91 | 7 | 30 | 10-178 | |
| o-Xylene | 22.6 | 21.5 | 95 | 10 | 30 | 29-134 | |
| sec-Butylbenzene | 22.6 | 20.6 | 91 | 6 | 30 | 10-172 | |
| Styrene | 22.6 | 19.5 | 86 | 9 | 30 | 27-127 | |
| Tetrachloroethene | 22.6 | 21.7 | 96 | 0 | 30 | 31-135 | |
| Toluene | 22.6 | 21.7 | 96 | 7 | 30 | 39-129 | |
| trans-1,2-Dichloroethene | 22.6 | 21.8 | 97 | 10 | 30 | 50-123 | |
| trans-1,3-Dichloropropene | 22.6 | 19.7 | 87 | 14 | 30 | 28-137 | |
| Trichloroethene | 22.6 | 22.5 | 99 | 10 | 30 | 10-177 | |
| Trichlorofluoromethane | 22.6 | 25.1 | 111 | 17 | 30 | 36-142 | |
| Vinyl acetate | 22.6 | 17.9 | 79 | 2 | 30 | 70-130 | |
| Vinyl chloride | 22.6 | 18.5 | 82 | 9 | 30 | 42-120 | |
| Xylenes, Total | 45.2 | 41.7 | 92 | 7 | 30 | 30-131 | |
| tert-Butylbenzene | 22.6 | 21.5 | 95 | 7 | 30 | 10-163 | |
| 1,2-Dibromoethane | 22.6 | 21.9 | 97 | 10 | 30 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 146331.D Lab Sample ID: MB 240-206089/1-A

Matrix: Solid Heated Purge:(Y/N) Y

Instrument ID: A3UX14 Date Analyzed: 11/10/2015 12:56

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 146336.D | 11/10/2015 14:43 |
| F-VB-UST-2-14 | 240-57597-2 | 146337.D | 11/10/2015 15:05 |
| F-VB-UST-3-14 | 240-57597-3 | 146338.D | 11/10/2015 15:26 |
| F-VB-UST-4-14 | 240-57597-4 | 146339.D | 11/10/2015 15:48 |
| F-VB-UST-5-14 | 240-57597-5 | 146340.D | 11/10/2015 16:09 |
| F-VB-UST-5-14 MS | 240-57597-5 MS | 146341.D | 11/10/2015 16:30 |
| F-VB-UST-5-14 MSD | 240-57597-5 MSD | 146342.D | 11/10/2015 16:52 |
| F-VB-UST-6-14 | 240-57597-6 | 146343.D | 11/10/2015 17:13 |
| F-VB-UST-7-14 | 240-57597-7 | 146344.D | 11/10/2015 17:35 |
| F-VB-UST-8-14 | 240-57597-8 | 146345.D | 11/10/2015 17:56 |
| F-VB-UST-9-14 | 240-57597-9 | 146346.D | 11/10/2015 18:18 |
| F-VB-UST-10-14 | 240-57597-10 | 146347.D | 11/10/2015 18:39 |
| F-VB-UST-11-14 | 240-57597-11 | 146348.D | 11/10/2015 19:01 |
| F-VB-UST-12-14 | 240-57597-12 | 146349.D | 11/10/2015 19:22 |
| F-VB-UST-13-14 | 240-57597-13 | 146350.D | 11/10/2015 19:44 |
| F-VB-UST-14-14 | 240-57597-14 | 146351.D | 11/10/2015 20:05 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206089/1-A

Matrix: Solid | Lab File ID: 146331.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5.0(g) | Date Analyzed: 11/10/2015 12:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.50 | U | 5.0 | 0.50 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.65 | U | 5.0 | 0.65 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.31 | U | 5.0 | 0.31 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.42 | U | 5.0 | 0.42 |
| 75-34-3 | 1,1-Dichloroethane | 0.34 | U | 5.0 | 0.34 |
| 75-35-4 | 1,1-Dichloroethene | 0.80 | U | 5.0 | 0.80 |
| 563-58-6 | 1,1-Dichloropropene | 0.37 | U | 5.0 | 0.37 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.50 | U | 5.0 | 0.50 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.78 | U | 5.0 | 0.78 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.46 | U | 5.0 | 0.46 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.38 | U | 5.0 | 0.38 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.58 | U | 5.0 | 0.58 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2.7 | U | 10 | 2.7 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.34 | U | 5.0 | 0.34 |
| 107-06-2 | 1,2-Dichloroethane | 0.46 | U | 5.0 | 0.46 |
| 78-87-5 | 1,2-Dichloropropane | 0.14 | U | 5.0 | 0.14 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.69 | U | 5.0 | 0.69 |
| 142-28-9 | 1,3-Dichloropropane | 0.40 | U | 5.0 | 0.40 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.28 | U | 5.0 | 0.28 |
| 594-20-7 | 2,2-Dichloropropane | 0.44 | U | 5.0 | 0.44 |
| 78-93-3 | 2-Butanone (MEK) | 2.83 | J | 20 | 1.1 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.64 | U | 50 | 0.64 |
| 95-49-8 | 2-Chlorotoluene | 0.30 | U | 5.0 | 0.30 |
| 591-78-6 | 2-Hexanone | 0.68 | U | 20 | 0.68 |
| 106-43-4 | 4-Chlorotoluene | 0.26 | U | 5.0 | 0.26 |
| 99-87-6 | p-Isopropyltoluene | 0.41 | U | 5.0 | 0.41 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 1.2 | U | 20 | 1.2 |
| 67-64-1 | Acetone | 5.7 | U | 20 | 5.7 |
| 71-43-2 | Benzene | 0.90 | U | 5.0 | 0.90 |
| 108-86-1 | Bromobenzene | 0.33 | U | 5.0 | 0.33 |
| 75-25-2 | Bromoform | 0.24 | U | 5.0 | 0.24 |
| 74-83-9 | Bromomethane | 0.42 | U | 5.0 | 0.42 |
| 75-15-0 | Carbon disulfide | 0.59 | U | 5.0 | 0.59 |
| 56-23-5 | Carbon tetrachloride | 0.65 | U | 5.0 | 0.65 |
| 108-90-7 | Chlorobenzene | 0.53 | U | 5.0 | 0.53 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206089/1-A

Matrix: Solid | Lab File ID: 146331.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5.0(g) | Date Analyzed: 11/10/2015 12:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.37 | U | 5.0 | 0.37 |
| 75-00-3 | Chloroethane | 0.44 | U | 5.0 | 0.44 |
| 67-66-3 | Chloroform | 0.36 | U | 5.0 | 0.36 |
| 74-87-3 | Chloromethane | 0.79 | U | 5.0 | 0.79 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.28 | U | 5.0 | 0.28 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.71 | U | 5.0 | 0.71 |
| 74-95-3 | Dibromomethane | 0.42 | U | 5.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.33 | U | 5.0 | 0.33 |
| 75-71-8 | Dichlorodifluoromethane | 0.28 | U | 5.0 | 0.28 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 5.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 5.7 | U | 200 | 5.7 |
| 100-41-4 | Ethylbenzene | 0.32 | U | 5.0 | 0.32 |
| 87-68-3 | Hexachlorobutadiene | 0.40 | U | 5.0 | 0.40 |
| 108-20-3 | Diisopropyl ether | 0.59 | U | 10 | 0.59 |
| 98-82-8 | Isopropylbenzene | 0.20 | U | 5.0 | 0.20 |
| 1634-04-4 | Methyl tert-butyl ether | 0.43 | U | 5.0 | 0.43 |
| 75-09-2 | Methylene Chloride | 0.74 | U | 5.0 | 0.74 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.42 | U | 10 | 0.42 |
| 91-20-3 | Naphthalene | 0.35 | U | 5.0 | 0.35 |
| 104-51-8 | n-Butylbenzene | 0.59 | U | 5.0 | 0.59 |
| 103-65-1 | N-Propylbenzene | 0.60 | U | 5.0 | 0.60 |
| 95-47-6 | o-Xylene | 0.57 | U | 5.0 | 0.57 |
| 135-98-8 | sec-Butylbenzene | 0.91 | U | 5.0 | 0.91 |
| 100-42-5 | Styrene | 0.40 | U | 5.0 | 0.40 |
| 994-05-8 | Tert-amyl methyl ether | 0.43 | U | 5.0 | 0.43 |
| 637-92-3 | Ethyl t-butyl ether | 0.41 | U | 5.0 | 0.41 |
| 127-18-4 | Tetrachloroethene | 0.80 | U | 5.0 | 0.80 |
| 108-88-3 | Toluene | 0.27 | U | 5.0 | 0.27 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.32 | U | 5.0 | 0.32 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.29 | U | 5.0 | 0.29 |
| 79-01-6 | Trichloroethene | 0.38 | U | 5.0 | 0.38 |
| 75-69-4 | Trichlorofluoromethane | 0.29 | U | 5.0 | 0.29 |
| 108-05-4 | Vinyl acetate | 0.66 | U | 10 | 0.66 |
| 75-01-4 | Vinyl chloride | 0.30 | U | 5.0 | 0.30 |
| 1330-20-7 | Xylenes, Total | 0.54 | U | 10 | 0.54 |
| 98-06-6 | tert-Butylbenzene | 0.67 | U | 5.0 | 0.67 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206089/1-A

Matrix: Solid | Lab File ID: 146331.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5.0(g) | Date Analyzed: 11/10/2015 12:56

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206051 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.53 | U | 5.0 | 0.53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 91 | | 67-125 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 88 | | 37-132 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 82 | | 52-136 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 91 | | 58-123 |

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146327.D

Lab ID: LCS 240-206051/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 20.0 | 20.4 | 102 | 80-120 | |
| 1,1,1-Trichloroethane | 20.0 | 22.0 | 110 | 77-126 | |
| 1,1,2,2-Tetrachloroethane | 20.0 | 19.6 | 98 | 77-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 20.0 | 20.9 | 105 | 80-138 | |
| 1,1-Dichloroethane | 20.0 | 18.5 | 93 | 76-120 | |
| 1,1-Dichloroethene | 20.0 | 19.1 | 95 | 75-135 | |
| 1,1-Dichloropropene | 20.0 | 20.3 | 102 | 80-120 | |
| 1,2,3-Trichlorobenzene | 20.0 | 19.5 | 98 | 61-121 | |
| 1,2,3-Trichloropropane | 20.0 | 20.3 | 102 | 80-126 | |
| 1,2,4-Trichlorobenzene | 20.0 | 20.6 | 103 | 64-124 | |
| 1,2,4-Trimethylbenzene | 20.0 | 20.3 | 101 | 80-129 | |
| 1,2-Dibromo-3-Chloropropane | 20.0 | 19.1 | 95 | 61-132 | |
| 1,2-Dichlorobenzene | 20.0 | 19.0 | 95 | 76-120 | |
| 1,2-Dichloroethane | 20.0 | 19.3 | 97 | 72-120 | |
| 1,2-Dichloropropane | 20.0 | 17.7 | 88 | 80-120 | |
| 1,3-Dichlorobenzene | 20.0 | 19.3 | 96 | 78-120 | |
| 1,3-Dichloropropane | 20.0 | 18.6 | 93 | 80-120 | |
| 1,4-Dichlorobenzene | 20.0 | 18.8 | 94 | 75-120 | |
| 2,2-Dichloropropane | 20.0 | 21.3 | 106 | 69-135 | |
| 2-Butanone (MEK) | 40.0 | 31.5 | 79 | 52-131 | |
| 2-Chloroethyl vinyl ether | 20.0 | 14.8 J | 74 | 28-173 | |
| 2-Chlorotoluene | 20.0 | 20.2 | 101 | 78-120 | |
| 2-Hexanone | 40.0 | 31.3 | 78 | 64-136 | |
| 4-Chlorotoluene | 20.0 | 20.0 | 100 | 79-120 | |
| p-Isopropyltoluene | 20.0 | 20.7 | 103 | 77-131 | |
| 4-Methyl-2-pentanone (MIBK) | 40.0 | 30.7 | 77 | 67-135 | |
| Acetone | 40.0 | 32.8 | 82 | 41-137 | |
| Benzene | 20.0 | 18.5 | 93 | 79-120 | |
| Bromobenzene | 20.0 | 21.3 | 106 | 80-120 | |
| Bromoform | 20.0 | 16.3 | 82 | 62-133 | |
| Bromomethane | 20.0 | 18.8 | 94 | 42-136 | |
| Carbon disulfide | 20.0 | 15.1 | 75 | 62-146 | |
| Carbon tetrachloride | 20.0 | 21.9 | 110 | 71-129 | |
| Chlorobenzene | 20.0 | 19.9 | 99 | 78-120 | |
| Dibromochloromethane | 20.0 | 18.2 | 91 | 72-127 | |
| Chloroethane | 20.0 | 18.7 | 93 | 58-120 | |
| Chloroform | 20.0 | 19.9 | 100 | 77-120 | |
| Chloromethane | 20.0 | 14.5 | 72 | 50-120 | |
| cis-1,2-Dichloroethene | 20.0 | 19.6 | 98 | 76-120 | |
| cis-1,3-Dichloropropene | 20.0 | 17.3 | 86 | 74-128 | |
| Dibromomethane | 20.0 | 19.1 | 96 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 146327.D

Lab ID: LCS 240-206051/5 Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 20.0 | 21.1 | 106 | 79-120 | |
| Dichlorodifluoromethane | 20.0 | 16.4 | 82 | 26-120 | |
| Bromodichloromethane | 20.0 | 18.1 | 90 | 80-122 | |
| tert-Butyl alcohol (TBA) | 200 | 138 J | 69 | 60-125 | |
| Ethylbenzene | 20.0 | 19.8 | 99 | 79-120 | |
| Hexachlorobutadiene | 20.0 | 21.4 | 107 | 54-131 | |
| Isopropylbenzene | 20.0 | 20.6 | 103 | 76-122 | |
| Methyl tert-butyl ether | 20.0 | 18.7 | 93 | 49-165 | |
| Methylene Chloride | 20.0 | 16.7 | 84 | 75-120 | |
| m-Xylene & p-Xylene | 20.0 | 19.8 | 99 | 80-120 | |
| Naphthalene | 20.0 | 18.4 | 92 | 65-123 | |
| n-Butylbenzene | 20.0 | 20.2 | 101 | 68-135 | |
| N-Propylbenzene | 20.0 | 19.8 | 99 | 80-129 | |
| o-Xylene | 20.0 | 20.7 | 103 | 80-120 | |
| sec-Butylbenzene | 20.0 | 19.9 | 100 | 74-129 | |
| Styrene | 20.0 | 18.7 | 93 | 80-120 | |
| Tetrachloroethene | 20.0 | 19.6 | 98 | 79-120 | |
| Toluene | 20.0 | 20.5 | 102 | 75-120 | |
| trans-1,2-Dichloroethene | 20.0 | 20.1 | 100 | 78-120 | |
| trans-1,3-Dichloropropene | 20.0 | 19.3 | 97 | 73-131 | |
| Trichloroethene | 20.0 | 21.1 | 105 | 79-120 | |
| Trichlorofluoromethane | 20.0 | 22.3 | 111 | 57-146 | |
| Vinyl acetate | 20.0 | 16.9 | 85 | 70-130 | |
| Vinyl chloride | 20.0 | 15.8 | 79 | 57-120 | |
| Xylenes, Total | 40.0 | 40.5 | 101 | 80-120 | |
| tert-Butylbenzene | 20.0 | 20.1 | 101 | 76-126 | |
| 1,2-Dibromoethane | 20.0 | 19.7 | 99 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 146341.D

Lab ID: 240-57597-5 MS  Client ID: F-VB-UST-5-14 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 22.8 | 0.58 U | 23.0 | 101 | 34-135 | |
| 1,1,1-Trichloroethane | 22.8 | 0.75 U | 24.1 | 106 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 22.8 | 0.36 U | 26.6 | 117 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 22.8 | 0.49 U | 23.5 | 103 | 50-147 | |
| 1,1-Dichloroethane | 22.8 | 0.39 U | 21.1 | 93 | 54-122 | |
| 1,1-Dichloroethene | 22.8 | 0.93 U | 19.8 | 87 | 49-157 | |
| 1,1-Dichloropropene | 22.8 | 0.43 U | 22.7 | 100 | 50-122 | |
| 1,2,3-Trichlorobenzene | 22.8 | 0.58 U | 14.1 | 62 | 10-120 | |
| 1,2,3-Trichloropropane | 22.8 | 0.90 U | 23.8 | 104 | 32-174 | |
| 1,2,4-Trichlorobenzene | 22.8 | 0.44 U | 14.2 | 62 | 10-120 | |
| 1,2,4-Trimethylbenzene | 22.8 | 0.67 U | 18.7 | 82 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 22.8 | 3.1 U | 22.2 | 97 | 10-153 | |
| 1,2-Dichlorobenzene | 22.8 | 0.39 U | 18.1 | 79 | 17-122 | |
| 1,2-Dichloroethane | 22.8 | 0.53 U | 22.0 | 97 | 49-123 | |
| 1,2-Dichloropropane | 22.8 | 0.16 U | 20.8 | 91 | 61-120 | |
| 1,3-Dichlorobenzene | 22.8 | 0.80 U | 18.3 | 80 | 16-126 | |
| 1,3-Dichloropropane | 22.8 | 0.46 U | 22.5 | 98 | 54-128 | |
| 1,4-Dichlorobenzene | 22.8 | 0.32 U | 17.8 | 78 | 15-121 | |
| 2,2-Dichloropropane | 22.8 | 0.51 U | 24.0 | 105 | 49-132 | |
| 2-Butanone (MEK) | 45.6 | 4.3 J | 36.6 | 71 | 30-143 | |
| 2-Chloroethyl vinyl ether | 22.8 | 0.74 U | 16.4 J | 72 | 10-162 | |
| 2-Chlorotoluene | 22.8 | 0.35 U | 20.8 | 91 | 11-162 | |
| 2-Hexanone | 45.6 | 0.79 U | 52.2 | 114 | 37-147 | |
| 4-Chlorotoluene | 22.8 | 0.30 U | 19.4 | 85 | 15-149 | |
| p-Isopropyltoluene | 22.8 | 0.47 U | 19.0 | 83 | 10-165 | |
| 4-Methyl-2-pentanone (MIBK) | 45.6 | 1.4 U | 150 | 328 | 43-147 | F1 |
| Acetone | 45.6 | 13 J | 47.7 | 77 | 24-140 | |
| Benzene | 22.8 | 1.0 U | 21.1 | 93 | 53-120 | |
| Bromobenzene | 22.8 | 0.38 U | 21.7 | 95 | 24-144 | |
| Bromoform | 22.8 | 0.28 U | 18.7 | 82 | 18-129 | |
| Bromomethane | 22.8 | 0.49 U | 19.8 | 87 | 33-130 | |
| Carbon disulfide | 22.8 | 0.68 U | 16.0 | 70 | 20-151 | |
| Carbon tetrachloride | 22.8 | 0.75 U | 24.0 | 105 | 32-137 | |
| Chlorobenzene | 22.8 | 0.61 U | 20.7 | 91 | 37-120 | |
| Dibromochloromethane | 22.8 | 0.43 U | 20.2 | 89 | 29-135 | |
| Chloroethane | 22.8 | 0.51 U | 21.5 | 94 | 45-120 | |
| Chloroform | 22.8 | 0.42 U | 22.9 | 100 | 53-120 | |
| Chloromethane | 22.8 | 0.92 U | 15.9 | 70 | 34-120 | |
| cis-1,2-Dichloroethene | 22.8 | 0.32 U | 22.1 | 97 | 50-120 | |
| cis-1,3-Dichloropropene | 22.8 | 0.82 U | 19.1 | 84 | 27-133 | |
| Dibromomethane | 22.8 | 0.49 U | 22.0 | 96 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 146341.D

Lab ID: 240-57597-5 MS  Client ID: F-VB-UST-5-14 MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Bromochloromethane | 22.8 | 0.38 U | 23.3 | 102 | 53-120 | |
| Dichlorodifluoromethane | 22.8 | 0.32 U | 18.3 | 80 | 17-120 | |
| Bromodichloromethane | 22.8 | 0.34 U | 20.0 | 88 | 35-132 | |
| tert-Butyl alcohol (TBA) | 228 | 6.6 U | 164 J | 72 | 60-125 | |
| Ethylbenzene | 22.8 | 0.37 U | 20.5 | 90 | 30-131 | |
| Hexachlorobutadiene | 22.8 | 0.46 U | 12.5 | 55 | 10-131 | |
| Isopropylbenzene | 22.8 | 0.23 U | 19.5 | 86 | 21-134 | |
| Methyl tert-butyl ether | 22.8 | 0.50 U | 20.9 | 92 | 51-157 | |
| Methylene Chloride | 22.8 | 0.86 U | 18.8 | 82 | 54-120 | |
| m-Xylene & p-Xylene | 22.8 | 0.49 U | 19.3 | 85 | 29-131 | |
| Naphthalene | 22.8 | 0.41 J | 17.8 | 76 | 10-124 | |
| n-Butylbenzene | 22.8 | 0.68 U | 16.6 | 73 | 10-148 | |
| N-Propylbenzene | 22.8 | 0.70 U | 19.7 | 86 | 10-178 | |
| o-Xylene | 22.8 | 0.66 U | 20.9 | 92 | 29-134 | |
| sec-Butylbenzene | 22.8 | 1.1 U | 18.2 | 80 | 10-172 | |
| Styrene | 22.8 | 0.46 U | 17.9 | 79 | 27-127 | |
| Tetrachloroethene | 22.8 | 0.93 U | 22.3 | 98 | 31-135 | |
| Toluene | 22.8 | 0.31 U | 22.0 | 96 | 39-129 | |
| trans-1,2-Dichloroethene | 22.8 | 0.37 U | 22.3 | 98 | 50-123 | |
| trans-1,3-Dichloropropene | 22.8 | 0.34 U | 21.1 | 93 | 28-137 | |
| Trichloroethene | 22.8 | 0.44 U | 23.2 | 101 | 10-177 | |
| Trichlorofluoromethane | 22.8 | 0.34 U | 23.0 | 101 | 36-142 | |
| Vinyl acetate | 22.8 | 0.76 U | 14.4 | 63 | 70-130 | F1 |
| Vinyl chloride | 22.8 | 0.35 U | 17.4 | 76 | 42-120 | |
| Xylenes, Total | 45.6 | 0.63 U | 40.2 | 88 | 30-131 | |
| tert-Butylbenzene | 22.8 | 0.78 U | 21.3 | 93 | 10-163 | |
| 1,2-Dibromoethane | 22.8 | 0.61 U | 24.2 | 106 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 146342.D

Lab ID: 240-57597-5 MSD  Client ID: F-VB-UST-5-14 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 23.0 | 24.0 | 104 | 4 | 30 | 34-135 | |
| 1,1,1-Trichloroethane | 23.0 | 25.7 | 112 | 7 | 30 | 51-128 | |
| 1,1,2,2-Tetrachloroethane | 23.0 | 28.0 | 121 | 5 | 30 | 16-179 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 23.0 | 25.6 | 111 | 9 | 30 | 50-147 | |
| 1,1-Dichloroethane | 23.0 | 22.5 | 98 | 6 | 30 | 54-122 | |
| 1,1-Dichloroethene | 23.0 | 22.0 | 95 | 10 | 30 | 49-157 | |
| 1,1-Dichloropropene | 23.0 | 25.1 | 109 | 10 | 30 | 50-122 | |
| 1,2,3-Trichlorobenzene | 23.0 | 17.9 | 78 | 24 | 30 | 10-120 | |
| 1,2,3-Trichloropropane | 23.0 | 28.9 | 125 | 19 | 30 | 32-174 | |
| 1,2,4-Trichlorobenzene | 23.0 | 19.9 | 86 | 33 | 30 | 10-120 | F2 |
| 1,2,4-Trimethylbenzene | 23.0 | 25.2 | 109 | 29 | 30 | 10-173 | |
| 1,2-Dibromo-3-Chloropropane | 23.0 | 26.9 | 117 | 19 | 30 | 10-153 | |
| 1,2-Dichlorobenzene | 23.0 | 22.2 | 96 | 20 | 30 | 17-122 | |
| 1,2-Dichloroethane | 23.0 | 24.1 | 105 | 9 | 30 | 49-123 | |
| 1,2-Dichloropropane | 23.0 | 21.9 | 95 | 5 | 30 | 61-120 | |
| 1,3-Dichlorobenzene | 23.0 | 22.6 | 98 | 21 | 30 | 16-126 | |
| 1,3-Dichloropropane | 23.0 | 23.2 | 101 | 3 | 30 | 54-128 | |
| 1,4-Dichlorobenzene | 23.0 | 22.3 | 97 | 22 | 30 | 15-121 | |
| 2,2-Dichloropropane | 23.0 | 24.5 | 107 | 2 | 30 | 49-132 | |
| 2-Butanone (MEK) | 46.1 | 39.3 | 76 | 7 | 30 | 30-143 | |
| 2-Chloroethyl vinyl ether | 23.0 | 17.2 J | 75 | 5 | 30 | 10-162 | |
| 2-Chlorotoluene | 23.0 | 25.6 | 111 | 21 | 30 | 11-162 | |
| 2-Hexanone | 46.1 | 52.6 | 114 | 1 | 30 | 37-147 | |
| 4-Chlorotoluene | 23.0 | 24.9 | 108 | 25 | 30 | 15-149 | |
| p-Isopropyltoluene | 23.0 | 26.0 | 113 | 31 | 30 | 10-165 | F2 |
| 4-Methyl-2-pentanone (MIBK) | 46.1 | 140 | 303 | 7 | 30 | 43-147 | F1 |
| Acetone | 46.1 | 53.3 | 88 | 11 | 30 | 24-140 | |
| Benzene | 23.0 | 23.0 | 100 | 8 | 30 | 53-120 | |
| Bromobenzene | 23.0 | 26.9 | 117 | 22 | 30 | 24-144 | |
| Bromoform | 23.0 | 19.9 | 86 | 7 | 30 | 18-129 | |
| Bromomethane | 23.0 | 21.9 | 95 | 10 | 30 | 33-130 | |
| Carbon disulfide | 23.0 | 17.6 | 76 | 10 | 30 | 20-151 | |
| Carbon tetrachloride | 23.0 | 26.5 | 115 | 10 | 30 | 32-137 | |
| Chlorobenzene | 23.0 | 23.7 | 103 | 14 | 30 | 37-120 | |
| Dibromochloromethane | 23.0 | 21.2 | 92 | 5 | 30 | 29-135 | |
| Chloroethane | 23.0 | 22.3 | 97 | 4 | 30 | 45-120 | |
| Chloroform | 23.0 | 25.1 | 109 | 9 | 30 | 53-120 | |
| Chloromethane | 23.0 | 17.1 | 74 | 8 | 30 | 34-120 | |
| cis-1,2-Dichloroethene | 23.0 | 24.0 | 104 | 8 | 30 | 50-120 | |
| cis-1,3-Dichloropropene | 23.0 | 21.0 | 91 | 9 | 30 | 27-133 | |
| Dibromomethane | 23.0 | 24.2 | 105 | 9 | 30 | 52-123 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 146342.D

Lab ID: 240-57597-5 MSD  Client ID: F-VB-UST-5-14 MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Bromochloromethane | 23.0 | 26.0 | 113 | 11 | 30 | 53-120 | |
| Dichlorodifluoromethane | 23.0 | 18.5 | 80 | 1 | 30 | 17-120 | |
| Bromodichloromethane | 23.0 | 22.9 | 99 | 13 | 30 | 35-132 | |
| tert-Butyl alcohol (TBA) | 230 | 189 J | 82 | 14 | 30 | 60-125 | |
| Ethylbenzene | 23.0 | 24.4 | 106 | 17 | 30 | 30-131 | |
| Hexachlorobutadiene | 23.0 | 22.3 | 97 | 57 | 30 | 10-131 | F2 |
| Isopropylbenzene | 23.0 | 25.1 | 109 | 25 | 30 | 21-134 | |
| Methyl tert-butyl ether | 23.0 | 22.3 | 97 | 6 | 30 | 51-157 | |
| Methylene Chloride | 23.0 | 20.2 | 88 | 7 | 30 | 54-120 | |
| m-Xylene & p-Xylene | 23.0 | 24.0 | 104 | 22 | 30 | 29-131 | |
| Naphthalene | 23.0 | 21.0 | 89 | 16 | 30 | 10-124 | |
| n-Butylbenzene | 23.0 | 25.0 | 109 | 40 | 30 | 10-148 | F2 |
| N-Propylbenzene | 23.0 | 25.5 | 111 | 26 | 30 | 10-178 | |
| o-Xylene | 23.0 | 25.1 | 109 | 18 | 30 | 29-134 | |
| sec-Butylbenzene | 23.0 | 25.1 | 109 | 32 | 30 | 10-172 | F2 |
| Styrene | 23.0 | 21.9 | 95 | 20 | 30 | 27-127 | |
| Tetrachloroethene | 23.0 | 24.7 | 107 | 10 | 30 | 31-135 | |
| Toluene | 23.0 | 24.4 | 106 | 11 | 30 | 39-129 | |
| trans-1,2-Dichloroethene | 23.0 | 24.7 | 107 | 10 | 30 | 50-123 | |
| trans-1,3-Dichloropropene | 23.0 | 22.9 | 99 | 8 | 30 | 28-137 | |
| Trichloroethene | 23.0 | 25.3 | 110 | 9 | 30 | 10-177 | |
| Trichlorofluoromethane | 23.0 | 26.2 | 114 | 13 | 30 | 36-142 | |
| Vinyl acetate | 23.0 | 17.8 | 77 | 21 | 30 | 70-130 | |
| Vinyl chloride | 23.0 | 18.6 | 81 | 7 | 30 | 42-120 | |
| Xylenes, Total | 46.1 | 49.1 | 107 | 20 | 30 | 30-131 | |
| tert-Butylbenzene | 23.0 | 26.7 | 116 | 22 | 30 | 10-163 | |
| 1,2-Dibromoethane | 23.0 | 25.0 | 108 | 3 | 30 | 45-127 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM IV
GC/MS VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: UXC4396.D Lab Sample ID: MB 240-206461/6

Matrix: Water Heated Purge:(Y/N) N

Instrument ID: A3UX15 Date Analyzed: 11/12/2015 12:11

GC Column: DB-624 ID: 0.18(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-206461/4 | UXC4393.D | 11/12/2015 11:03 |
| TB-110515 | 240-57597-16 | UXC4397.D | 11/12/2015 12:33 |
| | 240-57564-B-1 MS | UXC4417.D | 11/12/2015 20:05 |
| | 240-57564-B-1 MSD | UXC4418.D | 11/12/2015 20:27 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206461/6

Matrix: Water | Lab File ID: UXC4396.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:11

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.28 | U | 1.0 | 0.28 |
| 71-55-6 | 1,1,1-Trichloroethane | 0.44 | U | 1.0 | 0.44 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.22 | U | 1.0 | 0.22 |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.45 | U | 1.0 | 0.45 |
| 75-34-3 | 1,1-Dichloroethane | 0.30 | U | 1.0 | 0.30 |
| 75-35-4 | 1,1-Dichloroethene | 0.45 | U | 1.0 | 0.45 |
| 563-58-6 | 1,1-Dichloropropene | 0.42 | U | 1.0 | 0.42 |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.37 | U | 1.0 | 0.37 |
| 96-18-4 | 1,2,3-Trichloropropane | 0.44 | U | 1.0 | 0.44 |
| 526-73-8 | 1,2,3-Trimethylbenzene | 0.47 | U | 5.0 | 0.47 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.32 | U | 1.0 | 0.32 |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.41 | U | 1.0 | 0.41 |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.82 | U | 2.0 | 0.82 |
| 95-50-1 | 1,2-Dichlorobenzene | 0.25 | U | 1.0 | 0.25 |
| 107-06-2 | 1,2-Dichloroethane | 0.23 | U | 1.0 | 0.23 |
| 78-87-5 | 1,2-Dichloropropane | 0.25 | U | 1.0 | 0.25 |
| 541-73-1 | 1,3-Dichlorobenzene | 0.19 | U | 1.0 | 0.19 |
| 142-28-9 | 1,3-Dichloropropane | 0.19 | U | 1.0 | 0.19 |
| 106-46-7 | 1,4-Dichlorobenzene | 0.27 | U | 1.0 | 0.27 |
| 594-20-7 | 2,2-Dichloropropane | 0.26 | U | 1.0 | 0.26 |
| 78-93-3 | 2-Butanone (MEK) | 0.583 | J | 10 | 0.53 |
| 110-75-8 | 2-Chloroethyl vinyl ether | 0.59 | U | 10 | 0.59 |
| 95-49-8 | 2-Chlorotoluene | 0.40 | U | 1.0 | 0.40 |
| 591-78-6 | 2-Hexanone | 0.48 | U | 10 | 0.48 |
| 106-43-4 | 4-Chlorotoluene | 0.29 | U | 1.0 | 0.29 |
| 99-87-6 | p-Isopropyltoluene | 0.43 | U | 1.0 | 0.43 |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 0.99 | U | 10 | 0.99 |
| 67-64-1 | Acetone | 0.94 | U | 10 | 0.94 |
| 71-43-2 | Benzene | 0.35 | U | 1.0 | 0.35 |
| 108-86-1 | Bromobenzene | 0.35 | U | 1.0 | 0.35 |
| 75-25-2 | Bromoform | 0.56 | U | 1.0 | 0.56 |
| 74-83-9 | Bromomethane | 0.44 | U | 1.0 | 0.44 |
| 75-15-0 | Carbon disulfide | 0.38 | U | 1.0 | 0.38 |
| 56-23-5 | Carbon tetrachloride | 0.43 | U | 1.0 | 0.43 |
| 108-90-7 | Chlorobenzene | 0.25 | U | 1.0 | 0.25 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-206461/6

Matrix: Water | Lab File ID: UXC4396.D

Analysis Method: 8260C | Date Collected:

Sample wt/vol: 5(mL) | Date Analyzed: 11/12/2015 12:11

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: DB-624 ID: 0.18(mm)

% Moisture: | Level: (low/med) Low

Analysis Batch No.: 206461 | Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 124-48-1 | Dibromochloromethane | 0.43 | U | 1.0 | 0.43 |
| 75-00-3 | Chloroethane | 0.32 | U | 1.0 | 0.32 |
| 67-66-3 | Chloroform | 0.25 | U | 1.0 | 0.25 |
| 74-87-3 | Chloromethane | 0.44 | U | 1.0 | 0.44 |
| 156-59-2 | cis-1,2-Dichloroethene | 0.26 | U | 1.0 | 0.26 |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.46 | U | 1.0 | 0.46 |
| 74-95-3 | Dibromomethane | 0.42 | U | 1.0 | 0.42 |
| 74-97-5 | Bromochloromethane | 0.50 | U | 1.0 | 0.50 |
| 75-71-8 | Dichlorodifluoromethane | 0.32 | U | 1.0 | 0.32 |
| 75-27-4 | Bromodichloromethane | 0.29 | U | 1.0 | 0.29 |
| 75-65-0 | tert-Butyl alcohol (TBA) | 4.9 | U | 50 | 4.9 |
| 100-41-4 | Ethylbenzene | 0.25 | U | 1.0 | 0.25 |
| 87-68-3 | Hexachlorobutadiene | 0.35 | U | 1.0 | 0.35 |
| 108-20-3 | Diisopropyl ether | 0.50 | U | 10 | 0.50 |
| 98-82-8 | Isopropylbenzene | 0.35 | U | 1.0 | 0.35 |
| 1634-04-4 | Methyl tert-butyl ether | 0.20 | U | 1.0 | 0.20 |
| 75-09-2 | Methylene Chloride | 0.33 | U | 1.0 | 0.33 |
| 179601-23-1 | m-Xylene & p-Xylene | 0.24 | U | 2.0 | 0.24 |
| 91-20-3 | Naphthalene | 0.45 | U | 1.0 | 0.45 |
| 104-51-8 | n-Butylbenzene | 0.31 | U | 1.0 | 0.31 |
| 103-65-1 | N-Propylbenzene | 0.40 | U | 1.0 | 0.40 |
| 95-47-6 | o-Xylene | 0.25 | U | 1.0 | 0.25 |
| 135-98-8 | sec-Butylbenzene | 0.48 | U | 1.0 | 0.48 |
| 100-42-5 | Styrene | 0.45 | U | 1.0 | 0.45 |
| 994-05-8 | Tert-amyl methyl ether | 0.30 | U | 5.0 | 0.30 |
| 637-92-3 | Ethyl t-butyl ether | 0.23 | U | 5.0 | 0.23 |
| 127-18-4 | Tetrachloroethene | 0.31 | U | 1.0 | 0.31 |
| 108-88-3 | Toluene | 0.23 | U | 1.0 | 0.23 |
| 156-60-5 | trans-1,2-Dichloroethene | 0.30 | U | 1.0 | 0.30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.56 | U | 1.0 | 0.56 |
| 79-01-6 | Trichloroethene | 0.22 | U | 1.0 | 0.22 |
| 75-69-4 | Trichlorofluoromethane | 0.49 | U | 1.0 | 0.49 |
| 108-05-4 | Vinyl acetate | 0.41 | U | 2.0 | 0.41 |
| 75-01-4 | Vinyl chloride | 0.29 | U | 1.0 | 0.29 |
| 1330-20-7 | Xylenes, Total | 0.52 | U | 2.0 | 0.52 |
| 98-06-6 | tert-Butylbenzene | 0.41 | U | 1.0 | 0.41 |

FORM I
GC/MS VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-206461/6

Matrix: Water  Lab File ID: UXC4396.D

Analysis Method: 8260C  Date Collected:

Sample wt/vol: 5(mL)  Date Analyzed: 11/12/2015 12:11

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: DB-624  ID: 0.18(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206461  Units: ug/L

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 106-93-4 | 1,2-Dibromoethane | 0.32 | U | 1.0 | 0.32 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 2037-26-5 | Toluene-d8 (Surr) | 88 | | 80-120 |
| 1868-53-7 | Dibromofluoromethane (Surr) | 99 | | 79-120 |
| 460-00-4 | 4-Bromofluorobenzene (Surr) | 97 | | 61-120 |
| 17060-07-0 | 1,2-Dichloroethane-d4 (Surr) | 95 | | 78-125 |

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4393.D

Lab ID: LCS 240-206461/4 Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1,1,2-Tetrachloroethane | 10.0 | 11.8 | 118 | 80-120 | |
| 1,1,1-Trichloroethane | 10.0 | 11.1 | 111 | 77-123 | |
| 1,1,2,2-Tetrachloroethane | 10.0 | 9.38 | 94 | 71-123 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 10.0 | 9.57 | 96 | 67-138 | |
| 1,1-Dichloroethane | 10.0 | 9.43 | 94 | 79-125 | |
| 1,1-Dichloroethene | 10.0 | 8.73 | 87 | 76-124 | |
| 1,1-Dichloropropene | 10.0 | 10.4 | 104 | 80-128 | |
| 1,2,3-Trichlorobenzene | 10.0 | 8.64 | 86 | 66-120 | |
| 1,2,3-Trichloropropane | 10.0 | 11.3 | 113 | 72-126 | |
| 1,2,4-Trichlorobenzene | 10.0 | 8.66 | 87 | 61-120 | |
| 1,2,4-Trimethylbenzene | 10.0 | 9.93 | 99 | 76-120 | |
| 1,2-Dibromo-3-Chloropropane | 10.0 | 9.17 | 92 | 50-132 | |
| 1,2-Dichlorobenzene | 10.0 | 9.76 | 98 | 79-120 | |
| 1,2-Dichloroethane | 10.0 | 10.5 | 105 | 80-120 | |
| 1,2-Dichloropropane | 10.0 | 9.07 | 91 | 78-124 | |
| 1,3-Dichlorobenzene | 10.0 | 9.90 | 99 | 79-120 | |
| 1,3-Dichloropropane | 10.0 | 9.29 | 93 | 80-120 | |
| 1,4-Dichlorobenzene | 10.0 | 9.83 | 98 | 79-120 | |
| 2,2-Dichloropropane | 10.0 | 10.4 | 104 | 71-122 | |
| 2-Butanone (MEK) | 20.0 | 16.3 | 82 | 56-138 | |
| 2-Chloroethyl vinyl ether | 10.0 | 8.93 J | 89 | 54-139 | |
| 2-Chlorotoluene | 10.0 | 10.4 | 104 | 80-120 | |
| 2-Hexanone | 20.0 | 17.6 | 88 | 55-141 | |
| 4-Chlorotoluene | 10.0 | 9.89 | 99 | 80-120 | |
| p-Isopropyltoluene | 10.0 | 9.42 | 94 | 77-120 | |
| 4-Methyl-2-pentanone (MIBK) | 20.0 | 18.1 | 90 | 64-135 | |
| Acetone | 20.0 | 13.6 | 68 | 34-148 | |
| Benzene | 10.0 | 9.23 | 92 | 80-120 | |
| Bromobenzene | 10.0 | 10.7 | 107 | 80-120 | |
| Bromoform | 10.0 | 10.7 | 107 | 56-122 | |
| Bromomethane | 10.0 | 11.0 | 110 | 38-132 | |
| Carbon disulfide | 10.0 | 7.86 | 79 | 65-144 | |
| Carbon tetrachloride | 10.0 | 11.6 | 116 | 77-131 | |
| Chlorobenzene | 10.0 | 10.2 | 102 | 80-120 | |
| Dibromochloromethane | 10.0 | 10.6 | 106 | 74-120 | |
| Chloroethane | 10.0 | 9.73 | 97 | 36-126 | |
| Chloroform | 10.0 | 10.6 | 106 | 80-120 | |
| Chloromethane | 10.0 | 8.46 | 85 | 48-133 | |
| cis-1,2-Dichloroethene | 10.0 | 9.74 | 97 | 79-120 | |
| cis-1,3-Dichloropropene | 10.0 | 9.52 | 95 | 74-126 | |
| Dibromomethane | 10.0 | 10.7 | 107 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4393.D

Lab ID: LCS 240-206461/4 Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | LCS CONCENTRATION (ug/L) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Bromochloromethane | 10.0 | 10.8 | 108 | 80-120 | |
| Dichlorodifluoromethane | 10.0 | 10.5 | 105 | 23-136 | |
| Bromodichloromethane | 10.0 | 11.0 | 110 | 80-120 | |
| tert-Butyl alcohol (TBA) | 100 | 72.2 | 72 | 38-144 | |
| Ethylbenzene | 10.0 | 10.5 | 105 | 80-120 | |
| Hexachlorobutadiene | 10.0 | 8.37 | 84 | 62-120 | |
| Isopropylbenzene | 10.0 | 10.4 | 104 | 77-120 | |
| Methyl tert-butyl ether | 10.0 | 10.1 | 101 | 69-121 | |
| Methylene Chloride | 10.0 | 9.34 | 93 | 77-129 | |
| m-Xylene & p-Xylene | 10.0 | 10.0 | 100 | 80-120 | |
| Naphthalene | 10.0 | 8.44 | 84 | 53-121 | |
| n-Butylbenzene | 10.0 | 8.71 | 87 | 74-120 | |
| N-Propylbenzene | 10.0 | 10.2 | 102 | 79-123 | |
| o-Xylene | 10.0 | 10.4 | 104 | 80-120 | |
| sec-Butylbenzene | 10.0 | 9.27 | 93 | 77-120 | |
| Styrene | 10.0 | 9.85 | 99 | 76-122 | |
| Tetrachloroethene | 10.0 | 10.9 | 109 | 78-121 | |
| Toluene | 10.0 | 9.77 | 98 | 80-120 | |
| trans-1,2-Dichloroethene | 10.0 | 9.53 | 95 | 80-124 | |
| trans-1,3-Dichloropropene | 10.0 | 9.90 | 99 | 75-131 | |
| Trichloroethene | 10.0 | 11.1 | 111 | 80-121 | |
| Trichlorofluoromethane | 10.0 | 9.70 | 97 | 61-133 | |
| Vinyl acetate | 10.0 | 9.77 | 98 | 27-160 | |
| Vinyl chloride | 10.0 | 9.07 | 91 | 52-121 | |
| Xylenes, Total | 20.0 | 20.4 | 102 | 80-120 | |
| tert-Butylbenzene | 10.0 | 9.67 | 97 | 77-121 | |
| 1,2-Dibromoethane | 10.0 | 10.7 | 107 | 80-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

Sample is not in this SDG

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Water  Level: Low  Lab File ID: UXC4417.D

Lab ID: 240-57564-B-1 MS  Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 2500 | 110 U | 2570 | 103 | 69-122 | |
| 1,1,2,2-Tetrachloroethane | 2500 | 55 U | 2130 | 85 | 61-130 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 2500 | 110 U | 1850 | 74 | 44-140 | |
| 1,1-Dichloroethane | 2500 | 75 U | 2270 | 91 | 73-124 | |
| 1,1-Dichloroethene | 2500 | 110 U | 2020 | 81 | 67-124 | |
| 1,2,4-Trichlorobenzene | 2500 | 80 U | 2090 | 84 | 48-120 | |
| 1,2-Dibromo-3-Chloropropane | 2500 | 210 U | 2310 | 92 | 42-130 | |
| 1,2-Dichlorobenzene | 2500 | 63 U | 2300 | 92 | 67-118 | |
| 1,2-Dichloroethane | 2500 | 58 U | 2580 | 103 | 74-125 | |
| 1,2-Dichloropropane | 2500 | 63 U | 2170 | 87 | 73-122 | |
| 1,3-Dichlorobenzene | 2500 | 48 U | 2250 | 90 | 65-120 | |
| 1,4-Dichlorobenzene | 2500 | 68 U | 2290 | 92 | 66-120 | |
| 2-Butanone (MEK) | 5000 | 7000 | 12400 | 109 | 49-132 | |
| 2-Hexanone | 5000 | 120 U | 4460 | 89 | 49-142 | |
| 4-Methyl-2-pentanone (MIBK) | 5000 | 370 J | 5380 | 100 | 58-136 | |
| Acetone | 5000 | 9800 | 18700 | 179 | 32-126 | F1 |
| Benzene | 2500 | 88 U | 2220 | 89 | 73-121 | |
| Bromoform | 2500 | 140 U | 2500 | 100 | 45-121 | |
| Bromomethane | 2500 | 110 U | 2390 | 96 | 26-136 | |
| Carbon disulfide | 2500 | 95 U | 1870 | 75 | 54-144 | |
| Carbon tetrachloride | 2500 | 110 U | 2620 | 105 | 65-129 | |
| Chlorobenzene | 2500 | 63 U | 2390 | 96 | 72-120 | |
| Dibromochloromethane | 2500 | 110 U | 2460 | 98 | 62-122 | |
| Chloroethane | 2500 | 80 U | 2200 | 88 | 27-131 | |
| Chloroform | 2500 | 63 U | 2560 | 102 | 73-121 | |
| Chloromethane | 2500 | 110 U | 1980 | 79 | 39-134 | |
| cis-1,2-Dichloroethene | 2500 | 65 U | 2370 | 95 | 66-124 | |
| cis-1,3-Dichloropropene | 2500 | 120 U | 2250 | 90 | 60-120 | |
| Dichlorodifluoromethane | 2500 | 80 U | 1870 | 75 | 14-137 | |
| Bromodichloromethane | 2500 | 73 U | 2650 | 106 | 72-120 | |
| Ethylbenzene | 2500 | 63 U | 2440 | 97 | 68-121 | |
| Isopropylbenzene | 2500 | 88 U | 2380 | 95 | 61-122 | |
| Methyl tert-butyl ether | 2500 | 50 U | 2520 | 101 | 61-121 | |
| Methylene Chloride | 2500 | 83 U | 2350 | 94 | 70-124 | |
| m-Xylene & p-Xylene | 2500 | 150 J | 2540 | 96 | 66-123 | |
| o-Xylene | 2500 | 72 J | 2500 | 97 | 68-121 | |
| Styrene | 2500 | 110 U | 2320 | 93 | 64-126 | |
| Tetrachloroethene | 2500 | 78 U | 2450 | 98 | 59-125 | |
| Toluene | 2500 | 170 J | 2380 | 88 | 72-122 | |
| trans-1,2-Dichloroethene | 2500 | 75 U | 2340 | 94 | 72-125 | |
| trans-1,3-Dichloropropene | 2500 | 140 U | 2330 | 93 | 58-132 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4417.D

Lab ID: 240-57564-B-1 MS Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | SAMPLE CONCENTRATION (ug/L) | MS CONCENTRATION (ug/L) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Trichloroethene | 2500 | 55 U | 2580 | 103 | 61-129 | |
| Trichlorofluoromethane | 2500 | 120 U | 2020 | 81 | 49-133 | |
| Vinyl chloride | 2500 | 73 U | 2040 | 81 | 44-122 | |
| Xylenes, Total | 5000 | 220 J | 5040 | 96 | 67-122 | |
| 1,2-Dibromoethane | 2500 | 80 U | 2430 | 97 | 69-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Water  Level: Low  Lab File ID: UXC4418.D

Lab ID: 240-57564-B-1 MSD  Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1,1-Trichloroethane | 2500 | 2630 | 105 | 2 | 14 | 69-122 | |
| 1,1,2,2-Tetrachloroethane | 2500 | 2220 | 89 | 4 | 18 | 61-130 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 2500 | 2160 | 87 | 15 | 35 | 44-140 | |
| 1,1-Dichloroethane | 2500 | 2310 | 93 | 2 | 14 | 73-124 | |
| 1,1-Dichloroethene | 2500 | 2090 | 84 | 4 | 24 | 67-124 | |
| 1,2,4-Trichlorobenzene | 2500 | 2190 | 88 | 5 | 28 | 48-120 | |
| 1,2-Dibromo-3-Chloropropane | 2500 | 2510 | 100 | 8 | 24 | 42-130 | |
| 1,2-Dichlorobenzene | 2500 | 2370 | 95 | 3 | 15 | 67-118 | |
| 1,2-Dichloroethane | 2500 | 2590 | 104 | 0 | 24 | 74-125 | |
| 1,2-Dichloropropane | 2500 | 2330 | 93 | 7 | 15 | 73-122 | |
| 1,3-Dichlorobenzene | 2500 | 2270 | 91 | 1 | 15 | 65-120 | |
| 1,4-Dichlorobenzene | 2500 | 2330 | 93 | 2 | 16 | 66-120 | |
| 2-Butanone (MEK) | 5000 | 12800 | 116 | 3 | 19 | 49-132 | |
| 2-Hexanone | 5000 | 4330 | 87 | 3 | 27 | 49-142 | |
| 4-Methyl-2-pentanone (MIBK) | 5000 | 5460 | 102 | 2 | 32 | 58-136 | |
| Acetone | 5000 | 18900 | 182 | 1 | 28 | 32-126 | F1 |
| Benzene | 2500 | 2250 | 90 | 1 | 13 | 73-121 | |
| Bromoform | 2500 | 2490 | 99 | 1 | 19 | 45-121 | |
| Bromomethane | 2500 | 2400 | 96 | 0 | 35 | 26-136 | |
| Carbon disulfide | 2500 | 1920 | 77 | 3 | 34 | 54-144 | |
| Carbon tetrachloride | 2500 | 2790 | 111 | 6 | 20 | 65-129 | |
| Chlorobenzene | 2500 | 2400 | 96 | 1 | 15 | 72-120 | |
| Dibromochloromethane | 2500 | 2450 | 98 | 0 | 19 | 62-122 | |
| Chloroethane | 2500 | 2180 | 87 | 1 | 35 | 27-131 | |
| Chloroform | 2500 | 2580 | 103 | 1 | 17 | 73-121 | |
| Chloromethane | 2500 | 2060 | 82 | 4 | 20 | 39-134 | |
| cis-1,2-Dichloroethene | 2500 | 2420 | 97 | 2 | 22 | 66-124 | |
| cis-1,3-Dichloropropene | 2500 | 2330 | 93 | 4 | 21 | 60-120 | |
| Dichlorodifluoromethane | 2500 | 2230 | 89 | 18 | 34 | 14-137 | |
| Bromodichloromethane | 2500 | 2760 | 110 | 4 | 19 | 72-120 | |
| Ethylbenzene | 2500 | 2370 | 95 | 3 | 16 | 68-121 | |
| Isopropylbenzene | 2500 | 2400 | 96 | 1 | 20 | 61-122 | |
| Methyl tert-butyl ether | 2500 | 2570 | 103 | 2 | 12 | 61-121 | |
| Methylene Chloride | 2500 | 2330 | 93 | 1 | 14 | 70-124 | |
| m-Xylene & p-Xylene | 2500 | 2510 | 94 | 2 | 15 | 66-123 | |
| o-Xylene | 2500 | 2490 | 97 | 0 | 14 | 68-121 | |
| Styrene | 2500 | 2340 | 93 | 1 | 15 | 64-126 | |
| Tetrachloroethene | 2500 | 2490 | 100 | 2 | 20 | 59-125 | |
| Toluene | 2500 | 2430 | 91 | 2 | 15 | 72-122 | |
| trans-1,2-Dichloroethene | 2500 | 2400 | 96 | 3 | 25 | 72-125 | |
| trans-1,3-Dichloropropene | 2500 | 2310 | 92 | 1 | 22 | 58-132 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM III
GC/MS VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Water Level: Low Lab File ID: UXC4418.D

Lab ID: 240-57564-B-1 MSD Client ID:

| COMPOUND | SPIKE ADDED (ug/L) | MSD CONCENTRATION (ug/L) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Trichloroethene | 2500 | 2730 | 109 | 6 | 14 | 61-129 | |
| Trichlorofluoromethane | 2500 | 2240 | 90 | 10 | 25 | 49-133 | |
| Vinyl chloride | 2500 | 2050 | 82 | 1 | 35 | 44-122 | |
| Xylenes, Total | 5000 | 5000 | 96 | 1 | 14 | 67-122 | |
| 1,2-Dibromoethane | 2500 | 2470 | 99 | 2 | 24 | 69-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8260C

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB14045.D BFB Injection Date: 08/13/2015

Instrument ID: A3UX14 BFB Injection Time: 13:54

Analysis Batch No.: 193150

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 19.1 | |
| 75 | 30.0 - 60.0 % of mass 95 | 45.9 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.7 | |
| 173 | Less than 2.0 % of mass 174 | 0.0 | (0.1)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 70.4 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.1 | (7.3)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 67.7 | (96.1)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.5 | (6.6)2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | IC 240-193150/12 | 144576.D | 08/13/2015 | 17:50 |
| | IC 240-193150/13 | 144577.D | 08/13/2015 | 18:11 |
| | IC 240-193150/14 | 144578.D | 08/13/2015 | 18:33 |
| | IC 240-193150/15 | 144579.D | 08/13/2015 | 18:54 |
| | IC 240-193150/16 | 144580.D | 08/13/2015 | 19:16 |
| | IC 240-193150/17 | 144581.D | 08/13/2015 | 19:37 |
| | IC 240-193150/18 | 144582.D | 08/13/2015 | 19:59 |
| | IC 240-193150/19 | 144583.D | 08/13/2015 | 20:20 |
| | IC 240-193150/20 | 144584.D | 08/13/2015 | 20:42 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50  Calibration End Date: 08/13/2015 20:42  Calibration ID: 30158

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-193150/20 | 144584.D |
| Level 2 | IC 240-193150/19 | 144583.D |
| Level 3 | IC 240-193150/18 | 144582.D |
| Level 4 | IC 240-193150/17 | 144581.D |
| Level 5 | IC 240-193150/16 | 144580.D |
| Level 6 | IC 240-193150/15 | 144579.D |
| Level 7 | IC 240-193150/14 | 144578.D |
| Level 8 | IC 240-193150/13 | 144577.D |
| Level 9 | IC 240-193150/12 | 144576.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Acetonitrile | 0.0270<br>0.0217 | 0.0199<br>0.0226 | 0.0223<br>0.0219 | 0.0211<br>0.0222 | 0.0231 | Ave | | 0.0224 | | | | 8.8 | | 15.0 | | | |
| 2-Chloro-1,3-butadiene | 0.3946<br>0.4330 | 0.4036<br>0.4520 | 0.4843<br>0.4367 | 0.4330<br>0.4384 | 0.4314 | Ave | | 0.4341 | | | | 6.0 | | 15.0 | | | |
| Diisopropyl ether | 0.1991<br>0.2087 | 0.1920<br>0.2170 | 0.2189<br>0.2067 | 0.2161<br>0.2066 | 0.2062 | Ave | | 0.2079 | | | | 4.2 | | 15.0 | | | |
| Ethyl t-butyl ether | 0.8061<br>0.8145 | 0.8124<br>0.8609 | 0.8459<br>0.8423 | 0.8104<br>0.8475 | 0.7961 | Ave | | 0.8262 | | | | 2.8 | | 15.0 | | | |
| Ethyl acetate | 0.2082<br>0.1555 | 0.1520<br>0.1600 | 0.1502<br>0.1572 | 0.1518<br>0.1578 | 0.1525 | Ave | | 0.1606 | | | | 11.3 | | 15.0 | | | |
| Propionitrile | 0.0291<br>0.0267 | 0.0294<br>0.0281 | 0.0277<br>0.0272 | 0.0279<br>0.0272 | 0.0264 | Ave | | 0.0278 | | | | 3.7 | | 15.0 | | | |
| Methacrylonitrile | 0.1071<br>0.1082 | 0.1056<br>0.1106 | 0.1086<br>0.1043 | 0.1077<br>0.1034 | 0.1075 | Ave | | 0.1070 | | | | 2.1 | | 15.0 | | | |
| Tert-amyl methyl ether | 0.6617<br>0.6962 | 0.6767<br>0.7331 | 0.7307<br>0.7077 | 0.7151<br>0.7140 | 0.6826 | Ave | | 0.7020 | | | | 3.5 | | 15.0 | | | |
| n-Butanol | 0.0138<br>0.0096 | 0.0095<br>0.0106 | 0.0097<br>0.0104 | 0.0114<br>0.0105 | 0.0089 | Ave | | 0.0105 | | | | 13.7 | | 15.0 | | | |
| Ethyl acrylate | 0.2316<br>0.2216 | 0.2219<br>0.2309 | 0.2178<br>0.2294 | 0.2151<br>0.2318 | 0.2122 | Ave | | 0.2236 | | | | 3.4 | | 15.0 | | | |
| Methyl methacrylate | 0.1524<br>0.1436 | 0.1363<br>0.1478 | 0.1383<br>0.1452 | 0.1404<br>0.1466 | 0.1384 | Ave | | 0.1432 | | | | 3.7 | | 15.0 | | | |
| 2-Nitropropane | +++++<br>0.0226 | 0.0209<br>0.0266 | 0.0198<br>0.0269 | 0.0195<br>0.0236 | 0.0194 | Ave | | 0.0224 | | | | 13.7 | | 15.0 | | | |
| n-Butyl acetate | +++++<br>0.3716 | +++++<br>0.3944 | 0.4011<br>0.3902 | 0.3921<br>0.3867 | 0.3596 | Ave | | 0.3851 | | | | 3.7 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 193150

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 08/13/2015 17:50 Calibration End Date: 08/13/2015 20:42 Calibration ID: 30158

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1-Chlorohexane | 0.4929<br>0.4383 | 0.4320<br>0.4614 | 0.4729<br>0.4436 | 0.4373<br>0.4435 | 0.4315 | Ave | | 0.4504 | | | | 4.7 | | 15.0 | | | |
| Cyclohexanone | 0.0128<br>0.0108 | 0.0155<br>0.0119 | 0.0120<br>0.0114 | 0.0149<br>0.0114 | 0.0103 | Ave | | 0.0123 | | | | 14.4 | | 15.0 | | | |
| 1,2,3-Trimethylbenzene | 2.8189<br>3.0746 | 2.8992<br>3.2639 | 3.1500<br>3.1145 | 3.0040<br>3.0651 | 2.9168 | Ave | | 3.0341 | | | | 4.6 | | 15.0 | | | |
| 1,3,5-Trichlorobenzene | 1.1368<br>1.0908 | 1.0640<br>1.1383 | 1.1518<br>1.0486 | 1.1209<br>1.0391 | 1.0725 | Ave | | 1.0959 | | | | 3.9 | | 15.0 | | | |
| 2-Methylnaphthalene | 1.1258<br>1.2729 | 1.0980<br>1.4105 | 1.0984<br>1.2876 | 1.2519<br>1.2359 | 1.1966 | Ave | | 1.2197 | | | | 8.4 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB14070.D  BFB Injection Date: 09/02/2015

Instrument ID: A3UX14  BFB Injection Time: 13:49

Analysis Batch No.: 196026

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 21.6 | |
| 75 | 30.0 - 60.0 % of mass 95 | 45.0 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 7.0 | |
| 173 | Less than 2.0 % of mass 174 | 0.6 | (0.8)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 70.3 | |
| 175 | 5.0 - 9.0 % of mass 174 | 4.9 | (7.0)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 68.0 | (96.8)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.6 | (6.7)2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | IC 240-196026/3 | 144935.D | 09/02/2015 | 14:32 |
| | IC 240-196026/4 | 144936.D | 09/02/2015 | 14:54 |
| | IC 240-196026/5 | 144937.D | 09/02/2015 | 15:15 |
| | ICIS 240-196026/6 | 144938.D | 09/02/2015 | 15:37 |
| | IC 240-196026/7 | 144939.D | 09/02/2015 | 15:58 |
| | IC 240-196026/8 | 144940.D | 09/02/2015 | 16:20 |
| | IC 240-196026/9 | 144941.D | 09/02/2015 | 16:41 |
| | IC 240-196026/10 | 144942.D | 09/02/2015 | 17:03 |
| | IC 240-196026/11 | 144943.D | 09/02/2015 | 17:24 |
| | ICV 240-196026/13 | 144945.D | 09/02/2015 | 18:07 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | IC 240-196026/11 | 144943.D |
| Level 2 | IC 240-196026/10 | 144942.D |
| Level 3 | IC 240-196026/9 | 144941.D |
| Level 4 | IC 240-196026/8 | 144940.D |
| Level 5 | IC 240-196026/7 | 144939.D |
| Level 6 | ICIS 240-196026/6 | 144938.D |
| Level 7 | IC 240-196026/5 | 144937.D |
| Level 8 | IC 240-196026/4 | 144936.D |
| Level 9 | IC 240-196026/3 | 144935.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | +++++<br>0.2391 | +++++<br>0.2507 | 0.2470<br>0.2355 | 0.2306<br>0.2458 | 0.2468 | Ave | | 0.2422 | | | 0.1000 | 3.0 | | 15.0 | | | |
| Chloromethane | 0.3754<br>0.3823 | 0.4019<br>0.3850 | 0.3740<br>0.3707 | 0.3723<br>0.3714 | 0.3751 | Ave | | 0.3787 | | | 0.1000 | 2.6 | | 15.0 | | | |
| Vinyl chloride | 0.2535<br>0.2723 | 0.2872<br>0.2816 | 0.2779<br>0.2705 | 0.2715<br>0.2802 | 0.2773 | Ave | | 0.2747 | | | 0.1000 | 3.5 | | 15.0 | | | |
| Butadiene | 0.2421<br>0.3144 | 0.2898<br>0.3180 | 0.3006<br>0.3092 | 0.3016<br>0.3170 | 0.3102 | Ave | | 0.3003 | | | | 7.9 | | 15.0 | | | |
| Bromomethane | 0.1681<br>0.1234 | 0.1376<br>+++++ | 0.1308<br>+++++ | 0.1354<br>+++++ | 0.1348 | Ave | | 0.1383 | | | 0.0500 | 11.1 | | 15.0 | | | |
| Chloroethane | 0.1729<br>0.1392 | 0.1452<br>0.1270 | 0.1461<br>+++++ | 0.1422<br>+++++ | 0.1409 | Ave | | 0.1448 | | | | 9.6 | | 15.0 | | | |
| Dichlorofluoromethane | 0.3353<br>0.3214 | 0.3424<br>0.3238 | 0.3254<br>0.3013 | 0.3326<br>0.3083 | 0.3302 | Ave | | 0.3245 | | | | 4.0 | | 15.0 | | | |
| Trichlorofluoromethane | 0.2372<br>0.3085 | 0.2496<br>0.3130 | 0.2776<br>0.2931 | 0.2862<br>0.3038 | 0.2978 | Ave | | 0.2852 | | | 0.1000 | 9.2 | | 15.0 | | | |
| Ethyl ether | 0.2263<br>0.1633 | 0.1893<br>0.1738 | 0.1892<br>0.1620 | 0.1718<br>0.1640 | 0.1839 | Ave | | 0.1804 | | | | 11.2 | | 15.0 | | | |
| Acrolein | 0.0216<br>0.0171 | 0.0159<br>0.0175 | 0.0196<br>+++++ | 0.0166<br>+++++ | 0.0179 | Ave | | 0.0180 | | | | 10.9 | | 15.0 | | | |
| 1,1-Dichloroethene | 0.2079<br>0.2313 | 0.1781<br>0.2399 | 0.2268<br>0.2226 | 0.2292<br>0.2287 | 0.2305 | Ave | | 0.2217 | | | 0.1000 | 8.3 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.1337<br>0.1711 | 0.1398<br>0.1761 | 0.1720<br>0.1647 | 0.1739<br>0.1677 | 0.1631 | Ave | | 0.1625 | | | 0.0500 | 9.4 | | 15.0 | | | |
| Acetone | +++++<br>0.0712 | +++++<br>0.0711 | 0.1266<br>0.0622 | 0.0892<br>0.0631 | 0.0783 | Lin1 | 0.6353 | 0.0626 | | | 0.0500 | | | | 0.9980 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32 Calibration End Date: 09/02/2015 17:24 Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Iodomethane | 0.3031<br>0.3309 | 0.3161<br>0.3401 | 0.3156<br>0.3166 | 0.3107<br>0.3246 | 0.3168 | Ave | | 0.3194 | | | | 3.4 | | 15.0 | | | |
| Carbon disulfide | +++++<br>0.6938 | +++++<br>0.7470 | 0.5786<br>0.7111 | 0.6047<br>0.7368 | 0.6277 | Ave | | 0.6714 | | | 0.1000 | 10.0 | | 15.0 | | | |
| 3-Chloro-1-propene | +++++<br>0.1475 | +++++<br>0.1563 | 0.1275<br>0.1460 | 0.1372<br>0.1497 | 0.1368 | Ave | | 0.1430 | | | | 6.8 | | 15.0 | | | |
| Methyl acetate | 0.1615<br>0.1574 | 0.1594<br>0.1602 | 0.1513<br>0.1539 | 0.1584<br>0.1527 | 0.1554 | Ave | | 0.1567 | | | 0.1000 | 2.3 | | 15.0 | | | |
| Methylene Chloride | +++++<br>0.2588 | +++++<br>0.2535 | 0.5579<br>0.2285 | 0.3574<br>0.2329 | 0.3001 | Lin1 | 1.5559 | 0.2263 | | | 0.1000 | | | | 0.9990 | | 0.9900 |
| tert-Butyl alcohol (TBA) | 0.0223<br>0.0196 | 0.0192<br>0.0202 | 0.0186<br>0.0190 | 0.0196<br>0.0197 | 0.0203 | Ave | | 0.0198 | | | | 5.4 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.2358<br>0.2589 | 0.2329<br>0.2662 | 0.2511<br>0.2514 | 0.2599<br>0.2562 | 0.2510 | Ave | | 0.2515 | | | 0.1000 | 4.3 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.5407<br>0.5781 | 0.5287<br>0.5970 | 0.5676<br>0.5608 | 0.5729<br>0.5665 | 0.5729 | Ave | | 0.5650 | | | 0.1000 | 3.6 | | 15.0 | | | |
| Acrylonitrile | 0.0788<br>0.0761 | 0.0771<br>0.0770 | 0.0752<br>0.0728 | 0.0765<br>0.0724 | 0.0759 | Ave | | 0.0757 | | | | 2.7 | | 15.0 | | | |
| Hexane | +++++<br>0.0773 | +++++<br>0.0782 | 0.1040<br>0.0738 | 0.0811<br>0.0760 | 0.0806 | Ave | | 0.0816 | | | | 12.5 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.4672<br>0.5014 | 0.4509<br>0.5208 | 0.4998<br>0.4916 | 0.4999<br>0.5087 | 0.4920 | Ave | | 0.4925 | | | 0.1000 | 4.3 | | 15.0 | | | |
| Vinyl acetate | 0.0241<br>0.0308 | 0.0166<br>0.0307 | 0.0238<br>+++++ | 0.0268<br>+++++ | 0.0267 | Lin1 | -0.018 | 0.0303 | | | | | | | 0.9970 | | 0.9900 |
| 2,2-Dichloropropane | 0.1821<br>0.2315 | 0.1870<br>0.2370 | 0.2083<br>0.2070 | 0.2175<br>0.2121 | 0.2141 | Ave | | 0.2107 | | | | 8.5 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.2557<br>0.2665 | 0.2545<br>0.2811 | 0.2602<br>0.2612 | 0.2694<br>0.2684 | 0.2713 | Ave | | 0.2654 | | | 0.1000 | 3.2 | | 15.0 | | | |
| 2-Butanone (MEK) | +++++<br>0.0942 | 0.1118<br>0.0991 | 0.0952<br>0.0916 | 0.0894<br>0.0929 | 0.0959 | Ave | | 0.0962 | | | 0.0500 | 7.2 | | 15.0 | | | |
| Bromochloromethane | 0.0858<br>0.1003 | 0.0854<br>0.1062 | 0.0973<br>0.1000 | 0.0967<br>0.1004 | 0.0986 | Ave | | 0.0967 | | | | 7.1 | | 15.0 | | | |
| Tetrahydrofuran | +++++<br>0.0638 | 0.0727<br>0.0652 | 0.0687<br>0.0628 | 0.0716<br>0.0619 | 0.0652 | Ave | | 0.0665 | | | | 6.1 | | 15.0 | | | |
| Chloroform | 0.4109<br>0.3987 | 0.3844<br>0.4148 | 0.3916<br>0.3944 | 0.3857<br>0.4100 | 0.3904 | Ave | | 0.3979 | | | | 2.9 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.2815<br>0.3335 | 0.2455<br>0.3520 | 0.3014<br>0.3288 | 0.3164<br>0.3392 | 0.3164 | Ave | | 0.3127 | | | 0.1000 | 10.5 | | 15.0 | | | |
| Cyclohexane | 0.5387<br>0.6083 | 0.4961<br>0.6348 | 0.6113<br>0.5854 | 0.6257<br>0.6043 | 0.6056 | Ave | | 0.5900 | | | 0.1000 | 7.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32  Calibration End Date: 09/02/2015 17:24  Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Carbon tetrachloride | 0.1974<br>0.2765 | 0.2070<br>0.2949 | 0.2344<br>0.2806 | 0.2508<br>0.2887 | 0.2562 | Ave | | 0.2541 | | | 0.1000 | 13.8 | | 15.0 | | | |
| 1,1-Dichloropropene | +++++<br>0.3352 | +++++<br>0.3502 | 0.3354<br>0.3258 | 0.3276<br>0.3378 | 0.3200 | Ave | | 0.3331 | | | | 3.0 | | 15.0 | | | |
| Benzene | 0.9461<br>1.0076 | 0.9369<br>1.0637 | 0.9831<br>0.9992 | 1.0297<br>1.0364 | 0.9873 | Ave | | 0.9989 | | | 0.5000 | 4.1 | | 15.0 | | | |
| Isobutyl alcohol | 0.0079<br>0.0070 | 0.0073<br>0.0070 | 0.0070<br>0.0065 | 0.0074<br>0.0064 | 0.0078 | Ave | | 0.0071 | | | | 7.3 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.2922<br>0.2806 | 0.2538<br>0.2902 | 0.2809<br>0.2746 | 0.2857<br>0.2775 | 0.2775 | Ave | | 0.2792 | | | 0.1000 | 4.0 | | 15.0 | | | |
| n-Heptane | 0.0495<br>0.0719 | 0.0549<br>0.0771 | 0.0671<br>0.0732 | 0.0733<br>0.0762 | 0.0704 | Ave | | 0.0682 | | | | 14.1 | | 15.0 | | | |
| Trichloroethene | 0.2322<br>0.2352 | 0.1925<br>0.2515 | 0.2312<br>0.2339 | 0.2474<br>0.2400 | 0.2383 | Ave | | 0.2336 | | | 0.2000 | 7.2 | | 15.0 | | | |
| Methylcyclohexane | 0.3977<br>0.5007 | 0.4180<br>0.5258 | 0.5131<br>0.4892 | 0.4996<br>0.5023 | 0.4970 | Ave | | 0.4826 | | | 0.1000 | 9.1 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.2351<br>0.2464 | 0.2304<br>0.2703 | 0.2458<br>0.2552 | 0.2519<br>0.2638 | 0.2445 | Ave | | 0.2493 | | | 0.1000 | 5.1 | | 15.0 | | | |
| Dibromomethane | 0.1138<br>0.1049 | 0.0928<br>0.1099 | 0.1030<br>0.1045 | 0.0989<br>0.1052 | 0.1030 | Ave | | 0.1040 | | | | 5.7 | | 15.0 | | | |
| 1,4-Dioxane | 0.0016<br>0.0015 | 0.0018<br>0.0016 | 0.0017<br>0.0016 | 0.0016<br>0.0016 | 0.0018 | Ave | | 0.0017 | | | | 5.4 | | 15.0 | | | |
| Bromodichloromethane | +++++<br>0.2583 | +++++<br>0.2789 | 0.2043<br>0.2740 | 0.2214<br>0.2839 | 0.2350 | Ave | | 0.2508 | | | 0.1000 | 12.3 | | 15.0 | | | |
| 2-Chloroethyl vinyl ether | 0.0880<br>0.0981 | 0.0738<br>0.1047 | 0.0849<br>0.1021 | 0.0908<br>0.1040 | 0.0933 | Ave | | 0.0933 | | | | 10.9 | | 15.0 | | | |
| cis-1,3-Dichloropropene | +++++<br>0.3137 | +++++<br>0.3452 | 0.2529<br>0.3385 | 0.2764<br>0.3487 | 0.2874 | Ave | | 0.3090 | | | 0.1500 | 12.2 | | 15.0 | | | |
| 4-Methyl-2-pentanone (MIBK) | 0.3475<br>0.3097 | 0.3025<br>0.3146 | 0.3102<br>0.2962 | 0.3082<br>0.2899 | 0.3133 | Ave | | 0.3102 | | | 0.0500 | 5.2 | | 15.0 | | | |
| Toluene | 1.6889<br>1.6108 | 1.5730<br>1.6794 | 1.5789<br>1.5858 | 1.6326<br>1.6292 | 1.5722 | Ave | | 1.6167 | | | 0.4000 | 2.8 | | 15.0 | | | |
| trans-1,3-Dichloropropene | +++++<br>0.3786 | +++++<br>0.4056 | 0.2981<br>0.3958 | 0.3438<br>0.4021 | 0.3521 | Ave | | 0.3680 | | | 0.1000 | 10.6 | | 15.0 | | | |
| Ethyl methacrylate | 0.3411<br>0.2949 | 0.2728<br>0.3134 | 0.2921<br>0.2986 | 0.3041<br>0.3000 | 0.2937 | Ave | | 0.3012 | | | | 6.1 | | 15.0 | | | |
| 1,1,2-Trichloroethane | 0.2387<br>0.2409 | 0.2020<br>0.2502 | 0.2341<br>0.2296 | 0.2342<br>0.2336 | 0.2367 | Ave | | 0.2333 | | | 0.1000 | 5.6 | | 15.0 | | | |
| Tetrachloroethene | 0.2604<br>0.2924 | 0.2503<br>0.3024 | 0.2921<br>0.2817 | 0.2907<br>0.2908 | 0.2885 | Ave | | 0.2833 | | | 0.2000 | 6.0 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32  Calibration End Date: 09/02/2015 17:24  Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,3-Dichloropropane | 0.4985<br>0.4202 | 0.3942<br>0.4386 | 0.3932<br>0.4193 | 0.4331<br>0.4249 | 0.4155 | Ave | | 0.4264 | | | | 7.3 | | 15.0 | | | |
| 2-Hexanone | 0.2259<br>0.1948 | 0.2056<br>0.1966 | 0.2059<br>0.1806 | 0.2139<br>0.1792 | 0.2041 | Ave | | 0.2007 | | | 0.0500 | 7.5 | | 15.0 | | | |
| Dibromochloromethane | +++++<br>0.2379 | 0.1676<br>0.2584 | 0.1731<br>0.2529 | 0.1970<br>0.2554 | 0.2047 | Lin1 | -0.324 | 0.2543 | | | | | | | 0.9990 | | 0.9900 |
| 1,2-Dibromoethane | 0.1915<br>0.2254 | 0.1731<br>0.2352 | 0.2140<br>0.2225 | 0.2199<br>0.2231 | 0.2195 | Ave | | 0.2138 | | | | 9.0 | | 15.0 | | | |
| Chlorobenzene | 0.8957<br>0.9330 | 0.9144<br>0.9681 | 0.9220<br>0.9247 | 0.9499<br>0.9450 | 0.9159 | Ave | | 0.9298 | | | 0.3000 | 2.3 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.2424<br>0.3183 | 0.2461<br>0.3292 | 0.2688<br>0.3070 | 0.2915<br>0.3071 | 0.3010 | Ave | | 0.2902 | | | | 10.7 | | 15.0 | | | |
| Ethylbenzene | 0.4967<br>0.5429 | 0.5117<br>0.5557 | 0.5343<br>0.5186 | 0.5496<br>0.5291 | 0.5275 | Ave | | 0.5296 | | | | 3.5 | | 15.0 | | | |
| m-Xylene & p-Xylene | 0.6493<br>0.6610 | 0.5935<br>0.6804 | 0.6630<br>0.6340 | 0.6529<br>0.6481 | 0.6392 | Ave | | 0.6468 | | | | 3.7 | | 15.0 | | | |
| o-Xylene | 0.6484<br>0.6613 | 0.5588<br>0.6783 | 0.6201<br>0.6144 | 0.6581<br>0.6145 | 0.6575 | Ave | | 0.6346 | | | | 5.7 | | 15.0 | | | |
| Styrene | 0.9630<br>0.9968 | 0.8800<br>1.0308 | 0.8828<br>0.9707 | 0.9401<br>0.9819 | 0.9540 | Ave | | 0.9556 | | | 0.3000 | 5.2 | | 15.0 | | | |
| Bromoform | 0.0617<br>0.1189 | 0.0736<br>0.1321 | 0.0749<br>0.1315 | 0.0937<br>0.1318 | 0.0977 | Lin1 | -0.136 | 0.1302 | | | 0.1000 | | | | 0.9970 | | 0.9900 |
| Isopropylbenzene | 1.6615<br>1.8186 | 1.5426<br>1.8815 | 1.7312<br>1.7266 | 1.7956<br>1.7460 | 1.7709 | Ave | | 1.7416 | | | 0.1000 | 5.6 | | 15.0 | | | |
| Bromobenzene | 0.7018<br>0.7269 | 0.6199<br>0.7522 | 0.6741<br>0.7263 | 0.7025<br>0.7372 | 0.6965 | Ave | | 0.7042 | | | | 5.6 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 0.5544<br>0.6344 | 0.5565<br>0.6397 | 0.5491<br>0.6163 | 0.5930<br>0.5999 | 0.6184 | Ave | | 0.5958 | | | 0.3000 | 5.9 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.1905<br>0.1847 | 0.1628<br>0.1813 | 0.1632<br>0.1748 | 0.1767<br>0.1723 | 0.1805 | Ave | | 0.1763 | | | | 5.3 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | 0.2675<br>0.2192 | 0.1583<br>0.2222 | 0.1934<br>0.2263 | 0.2038<br>0.2273 | 0.2186 | Ave | | 0.2152 | | | | 13.7 | | 15.0 | | | |
| N-Propylbenzene | 0.9746<br>1.0339 | 0.8541<br>1.0565 | 0.8793<br>1.0140 | 0.9898<br>1.0407 | 0.9489 | Ave | | 0.9769 | | | | 7.3 | | 15.0 | | | |
| 2-Chlorotoluene | 0.8414<br>0.8319 | 0.6742<br>0.8400 | 0.7804<br>0.8160 | 0.8063<br>0.8183 | 0.8038 | Ave | | 0.8014 | | | | 6.4 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 2.7121<br>3.3648 | 2.7168<br>3.3834 | 2.9008<br>3.2276 | 3.1601<br>3.2885 | 3.1601 | Ave | | 3.1016 | | | | 8.4 | | 15.0 | | | |
| 4-Chlorotoluene | 0.8033<br>0.7833 | 0.7373<br>0.8244 | 0.7337<br>0.7938 | 0.7735<br>0.8011 | 0.7475 | Ave | | 0.7776 | | | | 4.1 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 196026

SDG No.:

Instrument ID: A3UX14  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) Y

Calibration Start Date: 09/02/2015 14:32  Calibration End Date: 09/02/2015 17:24  Calibration ID: 30511

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| tert-Butylbenzene | 2.3374<br>2.9631 | 2.4133<br>3.0556 | 2.6550<br>2.8376 | 2.8314<br>2.8634 | 2.8090 | Ave | | 2.7518 | | | | 8.7 | | 15.0 | | | |
| 1,2,4-Trimethylbenzene | 2.7336<br>3.2868 | 2.6205<br>3.3941 | 2.8875<br>3.2352 | 3.0975<br>3.2908 | 3.1820 | Ave | | 3.0809 | | | | 8.8 | | 15.0 | | | |
| sec-Butylbenzene | 3.5842<br>4.5688 | 3.5902<br>4.6258 | 3.9612<br>4.3978 | 4.2686<br>4.4271 | 4.3250 | Ave | | 4.1943 | | | | 9.4 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.4872<br>1.4944 | 1.4279<br>1.5057 | 1.4500<br>1.4639 | 1.4918<br>1.4763 | 1.4693 | Ave | | 1.4741 | | | 0.6000 | 1.7 | | 15.0 | | | |
| p-Isopropyltoluene | 2.6700<br>3.5326 | 2.7312<br>3.5868 | 3.1681<br>3.3752 | 3.3032<br>3.4312 | 3.3784 | Ave | | 3.2418 | | | | 10.2 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 1.6796<br>1.4771 | 1.4525<br>1.4817 | 1.3997<br>1.4447 | 1.4789<br>1.4779 | 1.4661 | Ave | | 1.4842 | | | 0.5000 | 5.2 | | 15.0 | | | |
| n-Butylbenzene | 2.6703<br>3.3398 | 2.7947<br>3.4180 | 3.0393<br>3.3135 | 3.1433<br>3.3757 | 3.2352 | Ave | | 3.1478 | | | | 8.4 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.5577<br>1.4312 | 1.4056<br>1.4532 | 1.3574<br>1.3836 | 1.4146<br>1.3837 | 1.4022 | Ave | | 1.4210 | | | 0.4000 | 4.1 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | 0.0691<br>0.0964 | 0.0727<br>0.1007 | 0.0775<br>0.1005 | 0.0817<br>0.0987 | 0.0866 | Ave | | 0.0871 | | | 0.0500 | 14.3 | | 15.0 | | | |
| 1,2,4-Trichlorobenzene | 0.9903<br>0.9952 | 0.9066<br>0.9801 | 0.9495<br>0.9392 | 1.0303<br>0.9542 | 1.0144 | Ave | | 0.9733 | | | 0.2000 | 4.0 | | 15.0 | | | |
| Hexachlorobutadiene | 0.3960<br>0.5853 | 0.4364<br>0.5796 | 0.5790<br>0.5455 | 0.5879<br>0.5625 | 0.5837 | Ave | | 0.5396 | | | | 13.3 | | 15.0 | | | |
| Naphthalene | 2.5487<br>2.3198 | 2.2039<br>2.2590 | 2.1592<br>2.1437 | 2.2755<br>2.0812 | 2.3749 | Ave | | 2.2629 | | | | 6.2 | | 15.0 | | | |
| 1,2,3-Trichlorobenzene | 0.9121<br>0.9513 | 0.9280<br>0.9117 | 0.9440<br>0.8603 | 0.9768<br>0.8520 | 0.9646 | Ave | | 0.9223 | | | | 4.7 | | 15.0 | | | |
| Dibromofluoromethane (Surr) | +++++<br>0.2047 | +++++<br>0.2121 | 0.2454<br>0.1982 | 0.2289<br>0.2019 | 0.2161 | Ave | | 0.2154 | | | | 7.8 | | 15.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | +++++<br>0.2276 | +++++<br>0.2392 | 0.3027<br>0.2282 | 0.2781<br>0.2330 | 0.2521 | Ave | | 0.2516 | | | | 11.4 | | 15.0 | | | |
| Toluene-d8 (Surr) | +++++<br>1.4443 | +++++<br>1.4774 | 1.8704<br>1.3884 | 1.6510<br>1.4120 | 1.5274 | Ave | | 1.5387 | | | | 11.1 | | 15.0 | | | |
| 4-Bromofluorobenzene (Surr) | +++++<br>1.0832 | +++++<br>1.0727 | 1.4341<br>1.0549 | 1.2044<br>1.0662 | 1.1169 | Ave | | 1.1475 | | | | 11.9 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-196026/13 Calibration Date: 09/02/2015 18:07

Instrument ID: A3UX14 Calib Start Date: 09/02/2015 14:32

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 09/02/2015 17:24

Lab File ID: 144945.D Conc. Units: ng/uL Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2422 | 0.2902 | 0.1000 | 0.0240 | 0.0200 | 19.8 | 20.0 |
| Chloromethane | Ave | 0.3787 | 0.4264 | 0.1000 | 0.0225 | 0.0200 | 12.6 | 20.0 |
| Vinyl chloride | Ave | 0.2747 | 0.2991 | 0.1000 | 0.0218 | 0.0200 | 8.9 | 20.0 |
| Butadiene | Ave | 0.3003 | 0.3280 | | 0.0218 | 0.0200 | 9.2 | 20.0 |
| Bromomethane | Ave | 0.1383 | 0.1509 | 0.0500 | 0.0218 | 0.0200 | 9.1 | 20.0 |
| Chloroethane | Ave | 0.1448 | 0.1548 | 0.0500 | 0.0214 | 0.0200 | 6.9 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3245 | 0.3543 | | 0.0218 | 0.0200 | 9.2 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2852 | 0.3259 | 0.1000 | 0.0229 | 0.0200 | 14.3 | 20.0 |
| Ethyl ether | Ave | 0.1804 | 0.2164 | | 0.0240 | 0.0200 | 19.9 | 20.0 |
| Acrolein | Ave | 0.0180 | 0.0235 | | 0.130 | 0.100 | 30.4* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2217 | 0.2411 | 0.1000 | 0.0218 | 0.0200 | 8.8 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1625 | 0.1744 | 0.0500 | 0.0215 | 0.0200 | 7.4 | 20.0 |
| Acetone | Lin1 | | 0.0867 | 0.0500 | 0.0453 | 0.0400 | 13.2 | 20.0 |
| Iodomethane | Ave | 0.3194 | 0.3613 | | 0.0226 | 0.0200 | 13.1 | 20.0 |
| Carbon disulfide | Ave | 0.6714 | 0.6482 | 0.1000 | 0.0193 | 0.0200 | -3.4 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1430 | 0.1394 | | 0.0195 | 0.0200 | -2.5 | 20.0 |
| Methyl acetate | Ave | 0.1567 | 0.1835 | 0.1000 | 0.117 | 0.100 | 17.1 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3235 | 0.1000 | 0.0217 | 0.0200 | 8.6 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0198 | 0.0221 | | 0.223 | 0.200 | 11.3 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2515 | 0.2789 | 0.1000 | 0.0222 | 0.0200 | 10.9 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.5650 | 0.6276 | 0.1000 | 0.0222 | 0.0200 | 11.1 | 20.0 |
| Acrylonitrile | Ave | 0.0757 | 0.0862 | | 0.228 | 0.200 | 13.8 | 20.0 |
| Hexane | Ave | 0.0816 | 0.0814 | | 0.0200 | 0.0200 | -0.1 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4925 | 0.5567 | 0.1000 | 0.0226 | 0.0200 | 13.0 | 20.0 |
| Vinyl acetate | Lin1 | | 0.0217 | | 0.0149 | 0.0200 | -25.6* | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2107 | 0.2260 | | 0.0214 | 0.0200 | 7.2 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2654 | 0.2924 | 0.1000 | 0.0220 | 0.0200 | 10.2 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0962 | 0.1043 | 0.0500 | 0.0433 | 0.0400 | 8.4 | 20.0 |
| Bromochloromethane | Ave | 0.0967 | 0.1063 | | 0.0220 | 0.0200 | 9.9 | 20.0 |
| Tetrahydrofuran | Ave | 0.0665 | 0.0705 | | 0.0424 | 0.0400 | 6.0 | 20.0 |
| Chloroform | Ave | 0.3979 | 0.4367 | 0.2000 | 0.0219 | 0.0200 | 9.7 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3127 | 0.3465 | 0.1000 | 0.0222 | 0.0200 | 10.8 | 20.0 |
| Cyclohexane | Ave | 0.5900 | 0.6590 | 0.1000 | 0.0223 | 0.0200 | 11.7 | 20.0 |
| Carbon tetrachloride | Ave | 0.2541 | 0.2901 | 0.1000 | 0.0228 | 0.0200 | 14.2 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3331 | 0.3884 | | 0.0233 | 0.0200 | 16.6 | 20.0 |
| Benzene | Ave | 0.999 | 1.113 | 0.5000 | 0.0223 | 0.0200 | 11.4 | 20.0 |
| Isobutyl alcohol | Ave | 0.0071 | 0.0082 | | 0.573 | 0.500 | 14.6 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.2792 | 0.3100 | 0.1000 | 0.0222 | 0.0200 | 11.0 | 20.0 |
| n-Heptane | Ave | 0.0682 | 0.0773 | | 0.0227 | 0.0200 | 13.4 | 20.0 |
| Trichloroethene | Ave | 0.2336 | 0.2694 | 0.2000 | 0.0231 | 0.0200 | 15.3 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-196026/13  Calibration Date: 09/02/2015 18:07

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 144945.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.4826 | 0.5370 | 0.1000 | 0.0223 | 0.0200 | 11.3 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2493 | 0.2803 | 0.1000 | 0.0225 | 0.0200 | 12.5 | 20.0 |
| Dibromomethane | Ave | 0.1040 | 0.1205 | | 0.0232 | 0.0200 | 15.9 | 20.0 |
| 1,4-Dioxane | Ave | 0.0017 | 0.0019 | | 0.462 | 0.400 | 15.4 | 20.0 |
| Bromodichloromethane | Ave | 0.2508 | 0.2588 | 0.1000 | 0.0206 | 0.0200 | 3.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0933 | 0.1063 | | 0.0228 | 0.0200 | 14.0 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3090 | 0.3296 | 0.1500 | 0.0213 | 0.0200 | 6.7 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3102 | 0.3490 | 0.0500 | 0.0450 | 0.0400 | 12.5 | 20.0 |
| Toluene | Ave | 1.617 | 1.748 | 0.4000 | 0.0216 | 0.0200 | 8.1 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3680 | 0.4131 | 0.1000 | 0.0224 | 0.0200 | 12.2 | 20.0 |
| Ethyl methacrylate | Ave | 0.3012 | 0.3245 | | 0.0215 | 0.0200 | 7.7 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2333 | 0.2665 | 0.1000 | 0.0228 | 0.0200 | 14.2 | 20.0 |
| Tetrachloroethene | Ave | 0.2833 | 0.3198 | 0.2000 | 0.0226 | 0.0200 | 12.9 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4264 | 0.4668 | | 0.0219 | 0.0200 | 9.5 | 20.0 |
| 2-Hexanone | Ave | 0.2007 | 0.2198 | 0.0500 | 0.0438 | 0.0400 | 9.5 | 20.0 |
| Dibromochloromethane | Lin1 | | 0.2290 | | 0.0193 | 0.0200 | -3.6 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2138 | 0.2523 | | 0.0236 | 0.0200 | 18.0 | 20.0 |
| Chlorobenzene | Ave | 0.9298 | 1.009 | 0.3000 | 0.0217 | 0.0200 | 8.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.2902 | 0.3344 | | 0.0231 | 0.0200 | 15.3 | 20.0 |
| Ethylbenzene | Ave | 0.5296 | 0.5798 | | 0.0219 | 0.0200 | 9.5 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6468 | 0.6849 | | 0.0212 | 0.0200 | 5.9 | 20.0 |
| o-Xylene | Ave | 0.6346 | 0.7209 | | 0.0227 | 0.0200 | 13.6 | 20.0 |
| Styrene | Ave | 0.9556 | 0.9735 | 0.3000 | 0.0204 | 0.0200 | 1.9 | 20.0 |
| Bromoform | Lin1 | | 0.1061 | 0.1000 | 0.0173 | 0.0200 | -13.4 | 20.0 |
| Isopropylbenzene | Ave | 1.742 | 1.946 | 0.1000 | 0.0223 | 0.0200 | 11.7 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.5958 | 0.6826 | 0.3000 | 0.0229 | 0.0200 | 14.6 | 20.0 |
| Bromobenzene | Ave | 0.7042 | 0.7858 | | 0.0223 | 0.0200 | 11.6 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1763 | 0.2024 | | 0.0230 | 0.0200 | 14.8 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2152 | 0.2320 | | 0.0216 | 0.0200 | 7.8 | 20.0 |
| N-Propylbenzene | Ave | 0.9769 | 1.052 | | 0.0215 | 0.0200 | 7.7 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8014 | 0.8578 | | 0.0214 | 0.0200 | 7.0 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.102 | 3.417 | | 0.0220 | 0.0200 | 10.2 | 20.0 |
| 4-Chlorotoluene | Ave | 0.7776 | 0.8277 | | 0.0213 | 0.0200 | 6.5 | 20.0 |
| tert-Butylbenzene | Ave | 2.752 | 3.074 | | 0.0223 | 0.0200 | 11.7 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.081 | 3.372 | | 0.0219 | 0.0200 | 9.5 | 20.0 |
| sec-Butylbenzene | Ave | 4.194 | 4.751 | | 0.0227 | 0.0200 | 13.3 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.474 | 1.583 | 0.6000 | 0.0215 | 0.0200 | 7.4 | 20.0 |
| p-Isopropyltoluene | Ave | 3.242 | 3.585 | | 0.0221 | 0.0200 | 10.6 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.484 | 1.542 | 0.5000 | 0.0208 | 0.0200 | 3.9 | 20.0 |
| n-Butylbenzene | Ave | 3.148 | 3.388 | | 0.0215 | 0.0200 | 7.6 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.421 | 1.521 | 0.4000 | 0.0214 | 0.0200 | 7.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-196026/13  Calibration Date: 09/02/2015 18:07

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 144945.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0871 | 0.0879 | 0.0500 | 0.0202 | 0.0200 | 0.9 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.9733 | 0.9859 | 0.2000 | 0.0203 | 0.0200 | 1.3 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5396 | 0.6222 | | 0.0231 | 0.0200 | 15.3 | 20.0 |
| Naphthalene | Ave | 2.263 | 2.304 | | 0.0204 | 0.0200 | 1.8 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.9223 | 0.9287 | | 0.0201 | 0.0200 | 0.7 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2154 | 0.2054 | | 0.0477 | 0.0500 | -4.6 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2516 | 0.2368 | | 0.0471 | 0.0500 | -5.9 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.539 | 1.437 | | 0.0467 | 0.0500 | -6.6 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.147 | 1.045 | | 0.0455 | 0.0500 | -8.9 | 20.0 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB14128.D BFB Injection Date: 11/10/2015

Instrument ID: A3UX14 BFB Injection Time: 10:04

Analysis Batch No.: 206051

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 19.2 | |
| 75 | 30.0 - 60.0 % of mass 95 | 45.9 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.7 | |
| 173 | Less than 2.0 % of mass 174 | 0.0 | (0.0)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 75.8 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.4 | (7.1)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 73.3 | (96.7)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.8 | (6.6)2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-206051/3 | 146325.D | 11/10/2015 | 10:47 |
| | CCV 240-206051/4 | 146326.D | 11/10/2015 | 11:08 |
| | LCS 240-206051/5 | 146327.D | 11/10/2015 | 11:30 |
| | MB 240-206089/1-A | 146331.D | 11/10/2015 | 12:56 |
| F-VB-UST-1-14 | 240-57597-1 | 146336.D | 11/10/2015 | 14:43 |
| F-VB-UST-2-14 | 240-57597-2 | 146337.D | 11/10/2015 | 15:05 |
| F-VB-UST-3-14 | 240-57597-3 | 146338.D | 11/10/2015 | 15:26 |
| F-VB-UST-4-14 | 240-57597-4 | 146339.D | 11/10/2015 | 15:48 |
| F-VB-UST-5-14 | 240-57597-5 | 146340.D | 11/10/2015 | 16:09 |
| F-VB-UST-5-14 MS | 240-57597-5 MS | 146341.D | 11/10/2015 | 16:30 |
| F-VB-UST-5-14 MSD | 240-57597-5 MSD | 146342.D | 11/10/2015 | 16:52 |
| F-VB-UST-6-14 | 240-57597-6 | 146343.D | 11/10/2015 | 17:13 |
| F-VB-UST-7-14 | 240-57597-7 | 146344.D | 11/10/2015 | 17:35 |
| F-VB-UST-8-14 | 240-57597-8 | 146345.D | 11/10/2015 | 17:56 |
| F-VB-UST-9-14 | 240-57597-9 | 146346.D | 11/10/2015 | 18:18 |
| F-VB-UST-10-14 | 240-57597-10 | 146347.D | 11/10/2015 | 18:39 |
| F-VB-UST-11-14 | 240-57597-11 | 146348.D | 11/10/2015 | 19:01 |
| F-VB-UST-12-14 | 240-57597-12 | 146349.D | 11/10/2015 | 19:22 |
| F-VB-UST-13-14 | 240-57597-13 | 146350.D | 11/10/2015 | 19:44 |
| F-VB-UST-14-14 | 240-57597-14 | 146351.D | 11/10/2015 | 20:05 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206051/3  Calibration Date: 11/10/2015 10:47

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146325.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2422 | 0.3028 | 0.1000 | 0.0250 | 0.0200 | 25.0* | 20.0 |
| Chloromethane | Ave | 0.3787 | 0.3469 | 0.1000 | 0.0183 | 0.0200 | -8.4 | 20.0 |
| Vinyl chloride | Ave | 0.2747 | 0.2618 | 0.1000 | 0.0191 | 0.0200 | -4.7 | 20.0 |
| Butadiene | Ave | 0.3003 | 0.2963 | | 0.0197 | 0.0200 | -1.3 | 20.0 |
| Bromomethane | Ave | 0.1383 | 0.1443 | 0.0500 | 0.0209 | 0.0200 | 4.3 | 20.0 |
| Chloroethane | Ave | 0.1448 | 0.1576 | 0.0500 | 0.0218 | 0.0200 | 8.9 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3245 | 0.3591 | | 0.0221 | 0.0200 | 10.7 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2852 | 0.3576 | 0.1000 | 0.0251 | 0.0200 | 25.4* | 20.0 |
| Ethyl ether | Ave | 0.1804 | 0.1677 | | 0.0465 | 0.0500 | -7.0 | 20.0 |
| Acrolein | Ave | 0.0180 | 0.0143 | | 0.199 | 0.250 | -20.5* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2217 | 0.2482 | 0.1000 | 0.0560 | 0.0500 | 12.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1625 | 0.1959 | 0.0500 | 0.0603 | 0.0500 | 20.6* | 20.0 |
| Acetone | Lin1 | | 0.0643 | 0.0500 | 0.0926 | 0.100 | -7.4 | 20.0 |
| Iodomethane | Ave | 0.3194 | 0.3367 | | 0.0527 | 0.0500 | 5.4 | 20.0 |
| Carbon disulfide | Ave | 0.6714 | 0.6672 | 0.1000 | 0.0497 | 0.0500 | -0.6 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1430 | 0.1423 | | 0.0497 | 0.0500 | -0.5 | 20.0 |
| Methyl acetate | Ave | 0.1567 | 0.1298 | 0.1000 | 0.207 | 0.250 | -17.2 | 20.0 |
| Methylene Chloride | Lin1 | | 0.2547 | 0.1000 | 0.0494 | 0.0500 | -1.2 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0198 | 0.0165 | | 0.415 | 0.500 | -17.1 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2515 | 0.2656 | 0.1000 | 0.0528 | 0.0500 | 5.6 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.5650 | 0.5586 | 0.1000 | 0.0494 | 0.0500 | -1.1 | 20.0 |
| Acrylonitrile | Ave | 0.0757 | 0.0653 | | 0.431 | 0.500 | -13.8 | 20.0 |
| Hexane | Ave | 0.0816 | 0.0754 | | 0.0462 | 0.0500 | -7.5 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4925 | 0.4861 | 0.1000 | 0.0493 | 0.0500 | -1.3 | 20.0 |
| Vinyl acetate | Lin1 | | 0.0338 | | 0.0562 | 0.0500 | 12.4 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2107 | 0.2815 | | 0.0668 | 0.0500 | 33.6* | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2654 | 0.2800 | 0.1000 | 0.0528 | 0.0500 | 5.5 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0962 | 0.0843 | 0.0500 | 0.0876 | 0.100 | -12.4 | 20.0 |
| Bromochloromethane | Ave | 0.0967 | 0.1095 | | 0.0566 | 0.0500 | 13.2 | 20.0 |
| Tetrahydrofuran | Ave | 0.0665 | 0.0555 | | 0.0835 | 0.100 | -16.5 | 20.0 |
| Chloroform | Ave | 0.3979 | 0.4268 | 0.2000 | 0.0536 | 0.0500 | 7.3 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3127 | 0.3856 | 0.1000 | 0.0616 | 0.0500 | 23.3* | 20.0 |
| Cyclohexane | Ave | 0.5900 | 0.5511 | 0.1000 | 0.0467 | 0.0500 | -6.6 | 20.0 |
| Carbon tetrachloride | Ave | 0.2541 | 0.3080 | 0.1000 | 0.0606 | 0.0500 | 21.2* | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3331 | 0.3407 | | 0.0511 | 0.0500 | 2.3 | 20.0 |
| Benzene | Ave | 0.999 | 1.002 | 0.5000 | 0.0502 | 0.0500 | 0.3 | 20.0 |
| Isobutyl alcohol | Ave | 0.0071 | 0.0058 | | 1.02 | 1.25 | -18.5 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.2792 | 0.2827 | 0.1000 | 0.0506 | 0.0500 | 1.2 | 20.0 |
| n-Heptane | Ave | 0.0682 | 0.0764 | | 0.0560 | 0.0500 | 12.1 | 20.0 |
| Trichloroethene | Ave | 0.2336 | 0.2579 | 0.2000 | 0.0552 | 0.0500 | 10.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206051/3  Calibration Date: 11/10/2015 10:47

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146325.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.4826 | 0.5136 | 0.1000 | 0.0532 | 0.0500 | 6.4 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2493 | 0.2442 | 0.1000 | 0.0490 | 0.0500 | -2.0 | 20.0 |
| Dibromomethane | Ave | 0.1040 | 0.1112 | | 0.0535 | 0.0500 | 6.9 | 20.0 |
| 1,4-Dioxane | Ave | 0.0017 | 0.0017 | | 1.03 | 1.00 | 3.3 | 20.0 |
| Bromodichloromethane | Ave | 0.2508 | 0.2740 | 0.1000 | 0.0546 | 0.0500 | 9.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0933 | 0.0821 | | 0.0880 | 0.100 | -12.0 | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3090 | 0.3112 | 0.1500 | 0.0504 | 0.0500 | 0.7 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3102 | 0.2676 | 0.0500 | 0.0863 | 0.100 | -13.7 | 20.0 |
| Toluene | Ave | 1.617 | 1.728 | 0.4000 | 0.0534 | 0.0500 | 6.9 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3680 | 0.3806 | 0.1000 | 0.0517 | 0.0500 | 3.4 | 20.0 |
| Ethyl methacrylate | Ave | 0.3012 | 0.2921 | | 0.0485 | 0.0500 | -3.0 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2333 | 0.2582 | 0.1000 | 0.0553 | 0.0500 | 10.6 | 20.0 |
| Tetrachloroethene | Ave | 0.2833 | 0.3043 | 0.2000 | 0.0537 | 0.0500 | 7.4 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4264 | 0.4318 | | 0.0506 | 0.0500 | 1.3 | 20.0 |
| 2-Hexanone | Ave | 0.2007 | 0.1769 | 0.0500 | 0.0881 | 0.100 | -11.9 | 20.0 |
| Dibromochloromethane | Lin1 | | 0.2579 | | 0.0520 | 0.0500 | 3.9 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2138 | 0.2359 | | 0.0552 | 0.0500 | 10.3 | 20.0 |
| Chlorobenzene | Ave | 0.9298 | 0.9905 | 0.3000 | 0.0533 | 0.0500 | 6.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.2902 | 0.3234 | | 0.0557 | 0.0500 | 11.5 | 20.0 |
| Ethylbenzene | Ave | 0.5296 | 0.5533 | | 0.0522 | 0.0500 | 4.5 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6468 | 0.6818 | | 0.0527 | 0.0500 | 5.4 | 20.0 |
| o-Xylene | Ave | 0.6346 | 0.6810 | | 0.0537 | 0.0500 | 7.3 | 20.0 |
| Styrene | Ave | 0.9556 | 1.032 | 0.3000 | 0.0540 | 0.0500 | 8.0 | 20.0 |
| Bromoform | Lin1 | | 0.1280 | 0.1000 | 0.0502 | 0.0500 | 0.3 | 20.0 |
| Isopropylbenzene | Ave | 1.742 | 1.907 | 0.1000 | 0.0547 | 0.0500 | 9.5 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.5958 | 0.6150 | 0.3000 | 0.0516 | 0.0500 | 3.2 | 20.0 |
| Bromobenzene | Ave | 0.7042 | 0.7856 | | 0.0558 | 0.0500 | 11.6 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1763 | 0.1912 | | 0.0542 | 0.0500 | 8.5 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2152 | 0.1899 | | 0.0441 | 0.0500 | -11.8 | 20.0 |
| N-Propylbenzene | Ave | 0.9769 | 1.061 | | 0.0543 | 0.0500 | 8.6 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8014 | 0.8759 | | 0.0546 | 0.0500 | 9.3 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.102 | 3.395 | | 0.0547 | 0.0500 | 9.5 | 20.0 |
| 4-Chlorotoluene | Ave | 0.7776 | 0.8372 | | 0.0538 | 0.0500 | 7.7 | 20.0 |
| tert-Butylbenzene | Ave | 2.752 | 3.033 | | 0.0551 | 0.0500 | 10.2 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.081 | 3.382 | | 0.0549 | 0.0500 | 9.8 | 20.0 |
| sec-Butylbenzene | Ave | 4.194 | 4.534 | | 0.0540 | 0.0500 | 8.1 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.474 | 1.485 | 0.6000 | 0.0504 | 0.0500 | 0.7 | 20.0 |
| p-Isopropyltoluene | Ave | 3.242 | 3.675 | | 0.0567 | 0.0500 | 13.4 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.484 | 1.475 | 0.5000 | 0.0497 | 0.0500 | -0.6 | 20.0 |
| n-Butylbenzene | Ave | 3.148 | 3.476 | | 0.0552 | 0.0500 | 10.4 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.421 | 1.407 | 0.4000 | 0.0495 | 0.0500 | -1.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206051/3  Calibration Date: 11/10/2015 10:47

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146325.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0871 | 0.0936 | 0.0500 | 0.0537 | 0.0500 | 7.4 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.9733 | 1.040 | 0.2000 | 0.0534 | 0.0500 | 6.8 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5396 | 0.6133 | | 0.0568 | 0.0500 | 13.7 | 20.0 |
| Naphthalene | Ave | 2.263 | 2.099 | | 0.0464 | 0.0500 | -7.2 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.9223 | 0.9394 | | 0.0509 | 0.0500 | 1.9 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2154 | 0.2257 | | 0.0524 | 0.0500 | 4.8 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2516 | 0.2552 | | 0.0507 | 0.0500 | 1.5 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.539 | 1.542 | | 0.0501 | 0.0500 | 0.2 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.147 | 1.068 | | 0.0465 | 0.0500 | -7.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206051/4  Calibration Date: 11/10/2015 11:08

Instrument ID: A3UX14  Calib Start Date: 08/13/2015 17:50

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 08/13/2015 20:42

Lab File ID: 146326.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0224 | 0.0216 | | 0.481 | 0.500 | -3.7 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.4341 | 0.4643 | | 0.0535 | 0.0500 | 7.0 | 20.0 |
| Diisopropyl ether | Ave | 0.2079 | 0.2164 | | 0.0520 | 0.0500 | 4.1 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8262 | 0.7658 | | 0.0463 | 0.0500 | -7.3 | 20.0 |
| Ethyl acetate | Ave | 0.1606 | 0.1374 | | 0.0855 | 0.100 | -14.5 | 20.0 |
| Propionitrile | Ave | 0.0278 | 0.0233 | | 0.420 | 0.500 | -16.0 | 20.0 |
| Methacrylonitrile | Ave | 0.1070 | 0.1039 | | 0.486 | 0.500 | -2.9 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.7020 | 0.6252 | | 0.0445 | 0.0500 | -10.9 | 20.0 |
| n-Butanol | Ave | 0.0105 | 0.0071 | | 0.846 | 1.25 | -32.4* | 20.0 |
| Methyl methacrylate | Ave | 0.1432 | 0.1329 | | 0.0928 | 0.100 | -7.2 | 20.0 |
| 2-Nitropropane | Ave | 0.0224 | 0.0235 | | 0.105 | 0.100 | 4.8 | 20.0 |
| 1-Chlorohexane | Ave | 0.4504 | 0.5066 | | 0.0562 | 0.0500 | 12.5 | 20.0 |
| Cyclohexanone | Ave | 0.0123 | 0.0082 | | 0.334 | 0.500 | -33.3* | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.034 | 3.484 | | 0.0574 | 0.0500 | 14.8 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.096 | 1.298 | | 0.0592 | 0.0500 | 18.4 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.220 | 1.397 | | 0.115 | 0.100 | 14.5 | 20.0 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB14129.D  BFB Injection Date: 11/11/2015

Instrument ID: A3UX14  BFB Injection Time: 09:51

Analysis Batch No.: 206267

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 17.9 | |
| 75 | 30.0 - 60.0 % of mass 95 | 44.2 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.3 | |
| 173 | Less than 2.0 % of mass 174 | 0.0 | (0.0)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 76.6 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.9 | (7.7)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 73.1 | (95.3)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.5 | (6.2)2 |

1-Value is % mass 174  2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-206267/3 | 146358.D | 11/11/2015 | 10:39 |
| | CCV 240-206267/4 | 146359.D | 11/11/2015 | 11:00 |
| | LCS 240-206267/5 | 146360.D | 11/11/2015 | 11:22 |
| | MB 240-206267/6 | 146361.D | 11/11/2015 | 11:43 |
| F-VB-UST-DUP | 240-57597-15 | 146365.D | 11/11/2015 | 13:09 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 146366.D | 11/11/2015 | 13:31 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 146367.D | 11/11/2015 | 13:52 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206267/3  Calibration Date: 11/11/2015 10:39

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146358.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.2422 | 0.2811 | 0.1000 | 0.0232 | 0.0200 | 16.0 | 20.0 |
| Chloromethane | Ave | 0.3787 | 0.3232 | 0.1000 | 0.0171 | 0.0200 | -14.6 | 20.0 |
| Vinyl chloride | Ave | 0.2747 | 0.2526 | 0.1000 | 0.0184 | 0.0200 | -8.0 | 20.0 |
| Butadiene | Ave | 0.3003 | 0.2846 | | 0.0190 | 0.0200 | -5.2 | 20.0 |
| Bromomethane | Ave | 0.1383 | 0.1381 | 0.0500 | 0.0200 | 0.0200 | -0.2 | 20.0 |
| Chloroethane | Ave | 0.1448 | 0.1525 | 0.0500 | 0.0211 | 0.0200 | 5.3 | 20.0 |
| Dichlorofluoromethane | Ave | 0.3245 | 0.3469 | | 0.0214 | 0.0200 | 6.9 | 20.0 |
| Trichlorofluoromethane | Ave | 0.2852 | 0.3317 | 0.1000 | 0.0233 | 0.0200 | 16.3 | 20.0 |
| Ethyl ether | Ave | 0.1804 | 0.1573 | | 0.0436 | 0.0500 | -12.8 | 20.0 |
| Acrolein | Ave | 0.0180 | 0.0124 | | 0.172 | 0.250 | -31.2* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.2217 | 0.2132 | 0.1000 | 0.0481 | 0.0500 | -3.8 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.1625 | 0.1832 | 0.0500 | 0.0564 | 0.0500 | 12.8 | 20.0 |
| Acetone | Lin1 | | 0.0558 | 0.0500 | 0.0790 | 0.100 | -21.0* | 20.0 |
| Iodomethane | Ave | 0.3194 | 0.3184 | | 0.0499 | 0.0500 | -0.3 | 20.0 |
| Carbon disulfide | Ave | 0.6714 | 0.6540 | 0.1000 | 0.0487 | 0.0500 | -2.6 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1430 | 0.1337 | | 0.0467 | 0.0500 | -6.5 | 20.0 |
| Methyl acetate | Ave | 0.1567 | 0.1142 | 0.1000 | 0.182 | 0.250 | -27.1* | 20.0 |
| Methylene Chloride | Lin1 | | 0.2445 | 0.1000 | 0.0472 | 0.0500 | -5.7 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0198 | 0.0143 | | 0.359 | 0.500 | -28.2* | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.2515 | 0.2557 | 0.1000 | 0.0508 | 0.0500 | 1.7 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.5650 | 0.5068 | 0.1000 | 0.0448 | 0.0500 | -10.3 | 20.0 |
| Acrylonitrile | Ave | 0.0757 | 0.0582 | | 0.384 | 0.500 | -23.1* | 20.0 |
| Hexane | Ave | 0.0816 | 0.0688 | | 0.0422 | 0.0500 | -15.6 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.4925 | 0.4603 | 0.1000 | 0.0467 | 0.0500 | -6.5 | 20.0 |
| Vinyl acetate | Lin1 | | 0.0268 | | 0.0448 | 0.0500 | -10.4 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.2107 | 0.2606 | | 0.0618 | 0.0500 | 23.7* | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.2654 | 0.2665 | 0.1000 | 0.0502 | 0.0500 | 0.4 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0962 | 0.0686 | 0.0500 | 0.0713 | 0.100 | -28.7* | 20.0 |
| Bromochloromethane | Ave | 0.0967 | 0.1012 | | 0.0523 | 0.0500 | 4.6 | 20.0 |
| Tetrahydrofuran | Ave | 0.0665 | 0.0474 | | 0.0713 | 0.100 | -28.7* | 20.0 |
| Chloroform | Ave | 0.3979 | 0.4075 | 0.2000 | 0.0512 | 0.0500 | 2.4 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.3127 | 0.3691 | 0.1000 | 0.0590 | 0.0500 | 18.0 | 20.0 |
| Cyclohexane | Ave | 0.5900 | 0.5303 | 0.1000 | 0.0449 | 0.0500 | -10.1 | 20.0 |
| Carbon tetrachloride | Ave | 0.2541 | 0.2972 | 0.1000 | 0.0585 | 0.0500 | 17.0 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.3331 | 0.3194 | | 0.0479 | 0.0500 | -4.1 | 20.0 |
| Benzene | Ave | 0.999 | 0.9482 | 0.5000 | 0.0475 | 0.0500 | -5.1 | 20.0 |
| Isobutyl alcohol | Ave | 0.0071 | 0.0050 | | 0.873 | 1.25 | -30.2* | 20.0 |
| 1,2-Dichloroethane | Ave | 0.2792 | 0.2670 | 0.1000 | 0.0478 | 0.0500 | -4.4 | 20.0 |
| n-Heptane | Ave | 0.0682 | 0.0690 | | 0.0506 | 0.0500 | 1.2 | 20.0 |
| Trichloroethene | Ave | 0.2336 | 0.2428 | 0.2000 | 0.0520 | 0.0500 | 4.0 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206267/3  Calibration Date: 11/11/2015 10:39

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146358.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.4826 | 0.4880 | 0.1000 | 0.0506 | 0.0500 | 1.1 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2493 | 0.2230 | 0.1000 | 0.0447 | 0.0500 | -10.5 | 20.0 |
| Dibromomethane | Ave | 0.1040 | 0.1015 | | 0.0488 | 0.0500 | -2.4 | 20.0 |
| 1,4-Dioxane | Ave | 0.0017 | 0.0013 | | 0.789 | 1.00 | -21.1* | 20.0 |
| Bromodichloromethane | Ave | 0.2508 | 0.2514 | 0.1000 | 0.0501 | 0.0500 | 0.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.0933 | 0.0673 | | 0.0721 | 0.100 | -27.9* | 20.0 |
| cis-1,3-Dichloropropene | Ave | 0.3090 | 0.2735 | 0.1500 | 0.0443 | 0.0500 | -11.5 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.3102 | 0.2327 | 0.0500 | 0.0750 | 0.100 | -25.0* | 20.0 |
| Toluene | Ave | 1.617 | 1.639 | 0.4000 | 0.0507 | 0.0500 | 1.4 | 20.0 |
| trans-1,3-Dichloropropene | Ave | 0.3680 | 0.3380 | 0.1000 | 0.0459 | 0.0500 | -8.2 | 20.0 |
| Ethyl methacrylate | Ave | 0.3012 | 0.2523 | | 0.0419 | 0.0500 | -16.2 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.2333 | 0.2322 | 0.1000 | 0.0498 | 0.0500 | -0.5 | 20.0 |
| Tetrachloroethene | Ave | 0.2833 | 0.2858 | 0.2000 | 0.0505 | 0.0500 | 0.9 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.4264 | 0.3916 | | 0.0459 | 0.0500 | -8.2 | 20.0 |
| 2-Hexanone | Ave | 0.2007 | 0.1471 | 0.0500 | 0.0733 | 0.100 | -26.7* | 20.0 |
| Dibromochloromethane | Lin1 | | 0.2357 | | 0.0476 | 0.0500 | -4.8 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.2138 | 0.2082 | | 0.0487 | 0.0500 | -2.6 | 20.0 |
| Chlorobenzene | Ave | 0.9298 | 0.9062 | 0.3000 | 0.0487 | 0.0500 | -2.5 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.2902 | 0.3043 | | 0.0524 | 0.0500 | 4.9 | 20.0 |
| Ethylbenzene | Ave | 0.5296 | 0.5106 | | 0.0482 | 0.0500 | -3.6 | 20.0 |
| m-Xylene & p-Xylene | Ave | 0.6468 | 0.6472 | | 0.0500 | 0.0500 | 0.0 | 20.0 |
| o-Xylene | Ave | 0.6346 | 0.6427 | | 0.0506 | 0.0500 | 1.3 | 20.0 |
| Styrene | Ave | 0.9556 | 0.9588 | 0.3000 | 0.0502 | 0.0500 | 0.3 | 20.0 |
| Bromoform | Lin1 | | 0.1101 | 0.1000 | 0.0433 | 0.0500 | -13.4 | 20.0 |
| Isopropylbenzene | Ave | 1.742 | 1.827 | 0.1000 | 0.0524 | 0.0500 | 4.9 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.5958 | 0.5674 | 0.3000 | 0.0476 | 0.0500 | -4.8 | 20.0 |
| Bromobenzene | Ave | 0.7042 | 0.7254 | | 0.0515 | 0.0500 | 3.0 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.1763 | 0.1674 | | 0.0475 | 0.0500 | -5.0 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.2152 | 0.1618 | | 0.0376 | 0.0500 | -24.8* | 20.0 |
| N-Propylbenzene | Ave | 0.9769 | 1.026 | | 0.0525 | 0.0500 | 5.0 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8014 | 0.8181 | | 0.0510 | 0.0500 | 2.1 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.102 | 3.275 | | 0.0528 | 0.0500 | 5.6 | 20.0 |
| 4-Chlorotoluene | Ave | 0.7776 | 0.7791 | | 0.0501 | 0.0500 | 0.2 | 20.0 |
| tert-Butylbenzene | Ave | 2.752 | 2.911 | | 0.0529 | 0.0500 | 5.8 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.081 | 3.208 | | 0.0521 | 0.0500 | 4.1 | 20.0 |
| sec-Butylbenzene | Ave | 4.194 | 4.345 | | 0.0518 | 0.0500 | 3.6 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.474 | 1.377 | 0.6000 | 0.0467 | 0.0500 | -6.6 | 20.0 |
| p-Isopropyltoluene | Ave | 3.242 | 3.491 | | 0.0538 | 0.0500 | 7.7 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.484 | 1.348 | 0.5000 | 0.0454 | 0.0500 | -9.2 | 20.0 |
| n-Butylbenzene | Ave | 3.148 | 3.277 | | 0.0520 | 0.0500 | 4.1 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.421 | 1.338 | 0.4000 | 0.0471 | 0.0500 | -5.8 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206267/3  Calibration Date: 11/11/2015 10:39

Instrument ID: A3UX14  Calib Start Date: 09/02/2015 14:32

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 09/02/2015 17:24

Lab File ID: 146358.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Ave | 0.0871 | 0.0827 | 0.0500 | 0.0475 | 0.0500 | -5.1 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.9733 | 0.9676 | 0.2000 | 0.0497 | 0.0500 | -0.6 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5396 | 0.5669 | | 0.0525 | 0.0500 | 5.1 | 20.0 |
| Naphthalene | Ave | 2.263 | 1.925 | | 0.0425 | 0.0500 | -14.9 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 0.9223 | 0.8905 | | 0.0483 | 0.0500 | -3.5 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.2154 | 0.2264 | | 0.0526 | 0.0500 | 5.1 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.2516 | 0.2478 | | 0.0493 | 0.0500 | -1.5 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.539 | 1.561 | | 0.0507 | 0.0500 | 1.5 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 1.147 | 1.030 | | 0.0449 | 0.0500 | -10.3 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206267/4  Calibration Date: 11/11/2015 11:00

Instrument ID: A3UX14  Calib Start Date: 08/13/2015 17:50

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 08/13/2015 20:42

Lab File ID: 146359.D  Conc. Units: ng/uL  Heated Purge: (Y/N) Y

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0224 | 0.0212 | | 0.473 | 0.500 | -5.3 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.4341 | 0.4417 | | 0.0509 | 0.0500 | 1.7 | 20.0 |
| Diisopropyl ether | Ave | 0.2079 | 0.2137 | | 0.0514 | 0.0500 | 2.8 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8262 | 0.7386 | | 0.0447 | 0.0500 | -10.6 | 20.0 |
| Ethyl acetate | Ave | 0.1606 | 0.1311 | | 0.0817 | 0.100 | -18.3 | 20.0 |
| Propionitrile | Ave | 0.0278 | 0.0225 | | 0.405 | 0.500 | -18.9 | 20.0 |
| Methacrylonitrile | Ave | 0.1070 | 0.1032 | | 0.482 | 0.500 | -3.5 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.7020 | 0.6097 | | 0.0434 | 0.0500 | -13.1 | 20.0 |
| n-Butanol | Ave | 0.0105 | 0.0072 | | 0.856 | 1.25 | -31.5* | 20.0 |
| Methyl methacrylate | Ave | 0.1432 | 0.1282 | | 0.0895 | 0.100 | -10.5 | 20.0 |
| 2-Nitropropane | Ave | 0.0224 | 0.0219 | | 0.0976 | 0.100 | -2.4 | 20.0 |
| 1-Chlorohexane | Ave | 0.4504 | 0.4749 | | 0.0527 | 0.0500 | 5.5 | 20.0 |
| Cyclohexanone | Ave | 0.0123 | 0.0087 | | 0.353 | 0.500 | -29.3* | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.034 | 3.431 | | 0.0565 | 0.0500 | 13.1 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.096 | 1.271 | | 0.0580 | 0.0500 | 16.0 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.220 | 1.395 | | 0.114 | 0.100 | 14.4 | 20.0 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB460.D BFB Injection Date: 10/28/2015

Instrument ID: A3UX15 BFB Injection Time: 12:37

Analysis Batch No.: 204067

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 17.7 | |
| 75 | 30.0 - 60.0 % of mass 95 | 53.2 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.4 | |
| 173 | Less than 2.0 % of mass 174 | 0.7 | (0.9)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 73.2 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.2 | (7.1)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 70.9 | (96.9)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 4.6 | (6.5)2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD8260 240-204067/2 | UXC3824A.D | 10/28/2015 | 13:05 |
| | STD8260 240-204067/3 | UXC3825A.D | 10/28/2015 | 13:28 |
| | STD8260 240-204067/4 | UXC3826A.D | 10/28/2015 | 13:51 |
| | STD8260 240-204067/5 | UXC3827A.D | 10/28/2015 | 14:13 |
| | STD8260 240-204067/6 | UXC3828A.D | 10/28/2015 | 14:36 |
| | STD8260 240-204067/7 | UXC3829A.D | 10/28/2015 | 14:59 |
| | ICV 240-204067/14 | UXC3830.D | 10/28/2015 | 15:21 |
| | STD 240-204067/8 | UXC3831.D | 10/28/2015 | 15:44 |
| | STD 240-204067/9 | UXC3832.D | 10/28/2015 | 16:06 |
| | STD 240-204067/10 | UXC3833.D | 10/28/2015 | 16:29 |
| | STD 240-204067/11 | UXC3834.D | 10/28/2015 | 16:52 |
| | STD 240-204067/12 | UXC3835.D | 10/28/2015 | 17:14 |
| | STD 240-204067/13 | UXC3836.D | 10/28/2015 | 17:37 |
| | ICV 240-204067/15 | UXC3837.D | 10/28/2015 | 18:00 |

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD8260 240-204067/7 | UXC3829A.D |
| Level 2 | STD8260 240-204067/6 | UXC3828A.D |
| Level 3 | STD8260 240-204067/5 | UXC3827A.D |
| Level 4 | STD8260 240-204067/4 | UXC3826A.D |
| Level 5 | STD8260 240-204067/3 | UXC3825A.D |
| Level 6 | STD8260 240-204067/2 | UXC3824A.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dichlorodifluoromethane | 0.3763 0.4630 | 0.3685 | 0.4347 | 0.4095 | 0.4415 | Ave | | 0.4156 | | | 0.1000 | 9.1 | | 15.0 | | | |
| Chloromethane | 0.3134 0.3006 | 0.2822 | 0.3097 | 0.2949 | 0.3057 | Ave | | 0.3011 | | | 0.1000 | 3.8 | | 15.0 | | | |
| Vinyl chloride | 0.3653 0.3315 | 0.3066 | 0.3371 | 0.3133 | 0.3248 | Ave | | 0.3298 | | | 0.1000 | 6.3 | | 15.0 | | | |
| Butadiene | 0.3149 0.2921 | 0.2799 | 0.2958 | 0.2812 | 0.2944 | Ave | | 0.2931 | | | | 4.3 | | 15.0 | | | |
| Bromomethane | 0.1768 0.1785 | 0.1856 | 0.1888 | 0.1677 | 0.1766 | Ave | | 0.1790 | | | 0.0500 | 4.2 | | 15.0 | | | |
| Chloroethane | 0.2170 0.1827 | 0.1810 | 0.1946 | 0.1792 | 0.1864 | Ave | | 0.1902 | | | | 7.5 | | 15.0 | | | |
| Dichlorofluoromethane | 0.5491 0.5336 | 0.5008 | 0.5423 | 0.5081 | 0.5307 | Ave | | 0.5274 | | | | 3.6 | | 15.0 | | | |
| Trichlorofluoromethane | 0.4290 0.5156 | 0.4293 | 0.4745 | 0.4580 | 0.4761 | Ave | | 0.4637 | | | 0.1000 | 7.1 | | 15.0 | | | |
| Ethyl ether | 0.2524 0.2058 | 0.2022 | 0.2174 | 0.2184 | 0.2166 | Ave | | 0.2188 | | | | 8.1 | | 15.0 | | | |
| Acrolein | 0.0194 0.0208 | 0.0195 | 0.0213 | 0.0216 | 0.0207 | Ave | | 0.0206 | | | | 4.5 | | 15.0 | | | |
| 1,1-Dichloroethene | 0.3192 0.3103 | 0.3026 | 0.3175 | 0.3087 | 0.3189 | Ave | | 0.3129 | | | 0.1000 | 2.2 | | 15.0 | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.2468 0.2655 | 0.2417 | 0.2611 | 0.2571 | 0.2662 | Ave | | 0.2564 | | | 0.0500 | 3.9 | | 15.0 | | | |
| Acetone | 0.1494 0.0756 | 0.1046 | 0.0806 | 0.0779 | 0.0760 | Lin1 | 0.1360 | 0.0728 | | | 0.0500 | | | | 0.9990 | | 0.9900 |
| Iodomethane | 0.4808 0.4866 | 0.4808 | 0.4981 | 0.4871 | 0.4948 | Ave | | 0.4880 | | | | 1.5 | | 15.0 | | | |
| Carbon disulfide | 0.8042 0.9355 | 0.7777 | 0.8325 | 0.8658 | 0.9300 | Ave | | 0.8576 | | | 0.1000 | 7.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 3-Chloro-1-propene | 0.1574<br>0.1826 | 0.1531 | 0.1725 | 0.1821 | 0.1872 | Ave | | 0.1725 | | | | 8.3 | | 15.0 | | | |
| Methyl acetate | 0.1529<br>0.1529 | 0.1449 | 0.1499 | 0.1485 | 0.1560 | Ave | | 0.1509 | | | 0.1000 | 2.6 | | 15.0 | | | |
| Methylene Chloride | 0.6600<br>0.3263 | 0.4591 | 0.4025 | 0.3526 | 0.3465 | Lin1 | 0.3272 | 0.3223 | | | 0.1000 | | | | 0.9990 | | 0.9900 |
| tert-Butyl alcohol (TBA) | 0.0126<br>0.0126 | 0.0137 | 0.0130 | 0.0136 | 0.0145 | Ave | | 0.0133 | | | | 5.5 | | 15.0 | | | |
| Methyl tert-butyl ether | 0.7083<br>0.8027 | 0.7349 | 0.7554 | 0.7782 | 0.8151 | Ave | | 0.7657 | | | 0.1000 | 5.3 | | 15.0 | | | |
| trans-1,2-Dichloroethene | 0.3546<br>0.3362 | 0.3350 | 0.3511 | 0.3434 | 0.3491 | Ave | | 0.3449 | | | 0.1000 | 2.3 | | 15.0 | | | |
| Acrylonitrile | 0.0807<br>0.0839 | 0.0777 | 0.0809 | 0.0821 | 0.0849 | Ave | | 0.0817 | | | | 3.1 | | 15.0 | | | |
| Hexane | 0.0897<br>0.0895 | 0.0775 | 0.0959 | 0.0883 | 0.0933 | Ave | | 0.0890 | | | | 7.1 | | 15.0 | | | |
| 1,1-Dichloroethane | 0.6004<br>0.5990 | 0.5656 | 0.6055 | 0.5988 | 0.6163 | Ave | | 0.5976 | | | 0.1000 | 2.9 | | 15.0 | | | |
| Vinyl acetate | 0.2891<br>0.3663 | 0.2933 | 0.3167 | 0.3246 | 0.3559 | Ave | | 0.3243 | | | | 9.8 | | 15.0 | | | |
| 2,2-Dichloropropane | 0.2966<br>0.3299 | 0.3036 | 0.3180 | 0.3379 | 0.3501 | Ave | | 0.3227 | | | | 6.4 | | 15.0 | | | |
| cis-1,2-Dichloroethene | 0.3967<br>0.3605 | 0.3546 | 0.3747 | 0.3635 | 0.3760 | Ave | | 0.3710 | | | 0.1000 | 4.1 | | 15.0 | | | |
| 2-Butanone (MEK) | 0.0890<br>0.0927 | 0.0889 | 0.0877 | 0.0928 | 0.0965 | Ave | | 0.0913 | | | 0.0500 | 3.7 | | 15.0 | | | |
| Bromochloromethane | 0.1697<br>0.1681 | 0.1526 | 0.1693 | 0.1646 | 0.1743 | Ave | | 0.1664 | | | | 4.5 | | 15.0 | | | |
| Tetrahydrofuran | 0.0551<br>0.0522 | 0.0497 | 0.0500 | 0.0514 | 0.0520 | Ave | | 0.0518 | | | | 3.7 | | 15.0 | | | |
| Chloroform | 0.6342<br>0.6461 | 0.6330 | 0.6483 | 0.6338 | 0.6635 | Ave | | 0.6432 | | | | 1.9 | | 15.0 | | | |
| 1,1,1-Trichloroethane | 0.4416<br>0.5089 | 0.4542 | 0.4774 | 0.5040 | 0.5219 | Ave | | 0.4847 | | | 0.1000 | 6.6 | | 15.0 | | | |
| Cyclohexane | 0.5510<br>0.5788 | 0.5254 | 0.5645 | 0.5688 | 0.5952 | Ave | | 0.5639 | | | 0.1000 | 4.3 | | 15.0 | | | |
| Carbon tetrachloride | 0.3590<br>0.4615 | 0.3535 | 0.4034 | 0.4258 | 0.4608 | Ave | | 0.4107 | | | 0.1000 | 11.6 | | 15.0 | | | |
| 1,1-Dichloropropene | 0.4475<br>0.4891 | 0.4505 | 0.4742 | 0.4878 | 0.4988 | Ave | | 0.4747 | | | | 4.5 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Isobutyl alcohol | 0.0045<br>0.0052 | 0.0043 | 0.0050 | 0.0051 | 0.0054 | Ave | | 0.0049 | | | | 8.3 | | 15.0 | | | |
| Benzene | 1.4285<br>1.4410 | 1.3752 | 1.3957 | 1.3924 | 1.4478 | Ave | | 1.4134 | | | 0.5000 | 2.1 | | 15.0 | | | |
| 1,2-Dichloroethane | 0.5642<br>0.5505 | 0.5164 | 0.5367 | 0.5384 | 0.5567 | Ave | | 0.5438 | | | 0.1000 | 3.1 | | 15.0 | | | |
| n-Heptane | 0.0744<br>0.0875 | 0.0820 | 0.0810 | 0.0857 | 0.0897 | Ave | | 0.0834 | | | | 6.6 | | 15.0 | | | |
| Trichloroethene | 0.3624<br>0.3674 | 0.3397 | 0.3792 | 0.3617 | 0.3763 | Ave | | 0.3644 | | | 0.2000 | 3.9 | | 15.0 | | | |
| Methylcyclohexane | 0.5343<br>0.5582 | 0.5050 | 0.5318 | 0.5510 | 0.5699 | Ave | | 0.5417 | | | 0.1000 | 4.3 | | 15.0 | | | |
| 1,2-Dichloropropane | 0.2971<br>0.3003 | 0.2855 | 0.3021 | 0.2932 | 0.3103 | Ave | | 0.2981 | | | 0.1000 | 2.8 | | 15.0 | | | |
| Dibromomethane | 0.1762<br>0.1826 | 0.1628 | 0.1757 | 0.1777 | 0.1863 | Ave | | 0.1769 | | | | 4.5 | | 15.0 | | | |
| 1,4-Dioxane | 0.0027<br>0.0022 | 0.0028 | 0.0024 | 0.0026 | 0.0028 | Ave | | 0.0026 | | | | 9.4 | | 15.0 | | | |
| Bromodichloromethane | 0.3493<br>0.4531 | 0.3322 | 0.3819 | 0.4022 | 0.4403 | Ave | | 0.3932 | | | 0.1000 | 12.3 | | 15.0 | | | |
| 2-Chloroethyl vinyl ether | 0.1365<br>0.1714 | 0.1411 | 0.1536 | 0.1565 | 0.1661 | Ave | | 0.1542 | | | | 8.8 | | 15.0 | | | |
| cis-1,3-Dichloropropene | 0.3364<br>0.4977 | 0.3567 | 0.4034 | 0.4277 | 0.4767 | Lin1 | -0.234 | 0.4896 | | | 0.1500 | | | | 0.9970 | | 0.9900 |
| 4-Methyl-2-pentanone (MIBK) | 0.1433<br>0.1852 | 0.1446 | 0.1677 | 0.1669 | 0.1802 | Ave | | 0.1647 | | | 0.0500 | 10.6 | | 15.0 | | | |
| Toluene | 1.9429<br>2.0302 | 1.8468 | 1.9605 | 1.9532 | 2.0824 | Ave | | 1.9693 | | | 0.4000 | 4.1 | | 15.0 | | | |
| trans-1,3-Dichloropropene | 0.3301<br>0.5505 | 0.3500 | 0.4184 | 0.4518 | 0.5264 | Lin1 | -0.328 | 0.5405 | | | 0.1000 | | | | 0.9960 | | 0.9900 |
| Ethyl methacrylate | 0.2971<br>0.4364 | 0.3159 | 0.3628 | 0.3705 | 0.4242 | Lin1 | -0.201 | 0.4307 | | | | | | | 0.9970 | | 0.9900 |
| 1,1,2-Trichloroethane | 0.3319<br>0.3389 | 0.3188 | 0.3304 | 0.3242 | 0.3461 | Ave | | 0.3317 | | | 0.1000 | 3.0 | | 15.0 | | | |
| Tetrachloroethene | 0.3571<br>0.3590 | 0.3460 | 0.3546 | 0.3549 | 0.3728 | Ave | | 0.3574 | | | 0.2000 | 2.5 | | 15.0 | | | |
| 1,3-Dichloropropane | 0.6595<br>0.6405 | 0.6138 | 0.6291 | 0.6236 | 0.6430 | Ave | | 0.6349 | | | | 2.5 | | 15.0 | | | |
| 2-Hexanone | 0.1368<br>0.1734 | 0.1326 | 0.1431 | 0.1652 | 0.1705 | Ave | | 0.1536 | | | 0.0500 | 11.8 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Dibromochloromethane | 0.2548<br>0.3871 | 0.2497 | 0.2924 | 0.3190 | 0.3650 | Lin1 | -0.213 | 0.3778 | | | | | | | 0.9950 | | 0.9900 |
| 1,2-Dibromoethane | 0.3173<br>0.3432 | 0.2888 | 0.3287 | 0.3283 | 0.3440 | Ave | | 0.3251 | | | | 6.3 | | 15.0 | | | |
| Chlorobenzene | 1.3001<br>1.3022 | 1.2131 | 1.2671 | 1.2315 | 1.2845 | Ave | | 1.2664 | | | 0.3000 | 2.9 | | 15.0 | | | |
| 1,1,1,2-Tetrachloroethane | 0.3166<br>0.4285 | 0.3268 | 0.3688 | 0.3781 | 0.4230 | Ave | | 0.3736 | | | | 12.5 | | 15.0 | | | |
| Ethylbenzene | 0.6139<br>0.6752 | 0.6272 | 0.6531 | 0.6630 | 0.6854 | Ave | | 0.6530 | | | | 4.2 | | 15.0 | | | |
| m-Xylene & p-Xylene | 1.6840<br>1.8801 | 1.6715 | 1.7805 | 1.7879 | 1.9286 | Ave | | 1.7888 | | | | 5.7 | | 15.0 | | | |
| o-Xylene | 0.7577<br>0.8160 | 0.7146 | 0.8018 | 0.7672 | 0.8323 | Ave | | 0.7816 | | | | 5.6 | | 15.0 | | | |
| Styrene | 1.2019<br>1.5179 | 1.2259 | 1.3763 | 1.3604 | 1.4960 | Ave | | 1.3631 | | | 0.3000 | 9.7 | | 15.0 | | | |
| Bromoform | 0.1273<br>0.1961 | 0.1074 | 0.1343 | 0.1474 | 0.1793 | Lin1 | -0.127 | 0.1882 | | | 0.1000 | | | | 0.9900 | | 0.9900 |
| Isopropylbenzene | 1.9913<br>2.0593 | 1.9528 | 2.1800 | 2.1276 | 2.3334 | Ave | | 2.1074 | | | 0.1000 | 6.6 | | 15.0 | | | |
| Bromobenzene | 0.9480<br>0.9172 | 0.8989 | 0.9318 | 0.9579 | 0.9573 | Ave | | 0.9352 | | | | 2.5 | | 15.0 | | | |
| 1,1,2,2-Tetrachloroethane | 0.6367<br>0.6917 | 0.6612 | 0.6532 | 0.6916 | 0.6953 | Ave | | 0.6716 | | | 0.3000 | 3.7 | | 15.0 | | | |
| 1,2,3-Trichloropropane | 0.2311<br>0.2455 | 0.2233 | 0.2502 | 0.2524 | 0.2593 | Ave | | 0.2436 | | | | 5.6 | | 15.0 | | | |
| trans-1,4-Dichloro-2-butene | 0.1847<br>0.2273 | 0.1712 | 0.1873 | 0.2097 | 0.2190 | Ave | | 0.1999 | | | | 11.0 | | 15.0 | | | |
| N-Propylbenzene | 0.9712<br>1.0332 | 0.9264 | 1.0226 | 1.0490 | 1.0539 | Ave | | 1.0094 | | | | 5.0 | | 15.0 | | | |
| 2-Chlorotoluene | 0.8756<br>0.8951 | 0.8629 | 0.9068 | 0.9045 | 0.9220 | Ave | | 0.8945 | | | | 2.4 | | 15.0 | | | |
| 1,3,5-Trimethylbenzene | 2.9650<br>3.6195 | 3.0996 | 3.3535 | 3.4351 | 3.5925 | Ave | | 3.3442 | | | | 7.9 | | 15.0 | | | |
| 4-Chlorotoluene | 3.4227<br>3.5911 | 3.2938 | 3.4666 | 3.4283 | 3.5772 | Ave | | 3.4633 | | | | 3.2 | | 15.0 | | | |
| tert-Butylbenzene | 2.6379<br>3.0255 | 2.5257 | 2.7206 | 2.8657 | 2.9264 | Ave | | 2.7836 | | | | 6.8 | | 15.0 | | | |
| 1,2,4-Trimethylbenzene | 3.2473<br>3.3932 | 3.1596 | 3.5475 | 3.6375 | 3.7459 | Ave | | 3.4552 | | | | 6.6 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 13:05 Calibration End Date: 10/28/2015 14:59 Calibration ID: 31159

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| sec-Butylbenzene | 3.5593<br>3.6545 | 3.5760 | 3.8326 | 3.9468 | 4.1043 | Ave | | 3.7789 | | | | 5.8 | | 15.0 | | | |
| 1,3-Dichlorobenzene | 1.9172<br>1.8494 | 1.7772 | 1.8392 | 1.8390 | 1.8648 | Ave | | 1.8478 | | | 0.6000 | 2.4 | | 15.0 | | | |
| p-Isopropyltoluene | 3.1956<br>3.3525 | 3.2534 | 3.4587 | 3.6196 | 3.7475 | Ave | | 3.4379 | | | | 6.2 | | 15.0 | | | |
| 1,4-Dichlorobenzene | 2.0031<br>1.8854 | 1.9004 | 1.8823 | 1.8816 | 1.9099 | Ave | | 1.9104 | | | 0.5000 | 2.4 | | 15.0 | | | |
| n-Butylbenzene | 2.6219<br>3.0543 | 2.5547 | 2.7935 | 2.9524 | 3.0366 | Ave | | 2.8356 | | | | 7.5 | | 15.0 | | | |
| 1,2-Dichlorobenzene | 1.8119<br>1.7466 | 1.7372 | 1.7233 | 1.7882 | 1.7748 | Ave | | 1.7637 | | | 0.4000 | 1.9 | | 15.0 | | | |
| 1,2-Dibromo-3-Chloropropane | 0.0810<br>0.1175 | 0.0782 | 0.0887 | 0.1005 | 0.1127 | Lin1 | -0.061 | 0.1154 | | | 0.0500 | | | | 0.9960 | | 0.9900 |
| 1,2,4-Trichlorobenzene | 1.2025<br>1.1327 | 1.0994 | 1.1644 | 1.1753 | 1.1939 | Ave | | 1.1614 | | | 0.2000 | 3.4 | | 15.0 | | | |
| Hexachlorobutadiene | 0.5980<br>0.4967 | 0.5848 | 0.5440 | 0.5646 | 0.5659 | Ave | | 0.5590 | | | | 6.4 | | 15.0 | | | |
| Naphthalene | 2.1201<br>2.7163 | 2.2108 | 2.4032 | 2.6263 | 2.8052 | Ave | | 2.4803 | | | | 11.3 | | 15.0 | | | |
| 1,2,3-Trichlorobenzene | 1.1193<br>0.9977 | 0.9897 | 1.0555 | 1.0704 | 1.0863 | Ave | | 1.0532 | | | | 4.8 | | 15.0 | | | |
| Dibromofluoromethane (Surr) | 0.3005<br>0.3229 | 0.3208 | 0.3302 | 0.3349 | 0.3342 | Ave | | 0.3239 | | | | 4.0 | | 15.0 | | | |
| 1,2-Dichloroethane-d4 (Surr) | 0.4565<br>0.4542 | 0.4715 | 0.4518 | 0.4649 | 0.4642 | Ave | | 0.4605 | | | | 1.6 | | 15.0 | | | |
| Toluene-d8 (Surr) | 1.7070<br>1.8254 | 1.7324 | 1.7232 | 1.7818 | 1.8396 | Ave | | 1.7682 | | | | 3.2 | | 15.0 | | | |
| 4-Bromofluorobenzene (Surr) | 0.6725<br>0.6778 | 0.6451 | 0.6646 | 0.6672 | 0.6928 | Ave | | 0.6700 | | | | 2.4 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 15:44 Calibration End Date: 10/28/2015 17:37 Calibration ID: 31161

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD 240-204067/13 | UXC3836.D |
| Level 2 | STD 240-204067/12 | UXC3835.D |
| Level 3 | STD 240-204067/11 | UXC3834.D |
| Level 4 | STD 240-204067/10 | UXC3833.D |
| Level 5 | STD 240-204067/9 | UXC3832.D |
| Level 6 | STD 240-204067/8 | UXC3831.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Acetonitrile | 0.0252 0.0219 | 0.0212 | 0.0233 | 0.0227 | 0.0226 | Ave | | 0.0228 | | | | 6.0 | | 15.0 | | | |
| Diisopropyl ether | 0.2683 0.2949 | 0.2607 | 0.2786 | 0.2899 | 0.2896 | Ave | | 0.2803 | | | | 4.8 | | 15.0 | | | |
| 2-Chloro-1,3-butadiene | 0.5072 0.5586 | 0.4581 | 0.5204 | 0.5317 | 0.5354 | Ave | | 0.5186 | | | | 6.6 | | 15.0 | | | |
| Ethyl t-butyl ether | 0.7639 0.8817 | 0.7574 | 0.8038 | 0.8303 | 0.8343 | Ave | | 0.8119 | | | | 5.8 | | 15.0 | | | |
| Ethyl acetate | 0.1723 0.1688 | 0.1526 | 0.1659 | 0.1628 | 0.1640 | Ave | | 0.1644 | | | | 4.1 | | 15.0 | | | |
| Propionitrile | 0.0300 0.0295 | 0.0271 | 0.0298 | 0.0298 | 0.0295 | Ave | | 0.0293 | | | | 3.7 | | 15.0 | | | |
| Methacrylonitrile | 0.1341 0.1392 | 0.1230 | 0.1302 | 0.1352 | 0.1327 | Ave | | 0.1324 | | | | 4.2 | | 15.0 | | | |
| Tert-amyl methyl ether | 0.6054 0.7084 | 0.5887 | 0.6367 | 0.6663 | 0.6719 | Ave | | 0.6462 | | | | 6.9 | | 15.0 | | | |
| n-Butanol | 0.0033 0.0038 | 0.0033 | 0.0034 | 0.0036 | 0.0038 | Ave | | 0.0035 | | | | 6.7 | | 15.0 | | | |
| Ethyl acrylate | 0.2445 0.2717 | 0.2373 | 0.2318 | 0.2579 | 0.2659 | Ave | | 0.2515 | | | | 6.4 | | 15.0 | | | |
| Methyl methacrylate | 0.1740 0.2003 | 0.1580 | 0.1764 | 0.1847 | 0.1885 | Ave | | 0.1803 | | | | 8.0 | | 15.0 | | | |
| 2-Nitropropane | 0.0283 0.0326 | 0.0236 | 0.0254 | 0.0265 | 0.0282 | Ave | | 0.0274 | | | | 11.3 | | 15.0 | | | |
| n-Butyl acetate | 0.2122 0.2685 | 0.2003 | 0.2057 | 0.2348 | 0.2498 | Ave | | 0.2286 | | | | 11.8 | | 15.0 | | | |
| 1-Chlorohexane | 0.4981 0.5513 | 0.4815 | 0.5004 | 0.5112 | 0.5405 | Ave | | 0.5138 | | | | 5.2 | | 15.0 | | | |
| Cyclohexanone | 0.0101 0.0133 | 0.0128 | 0.0128 | 0.0134 | 0.0144 | Ave | | 0.0128 | | | | 11.4 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204067

SDG No.:

Instrument ID: A3UX15 GC Column: DB-624 ID: 0.18(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/28/2015 15:44 Calibration End Date: 10/28/2015 17:37 Calibration ID: 31161

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,2,3-Trimethylbenzene | 3.2126<br>3.7399 | 3.1433 | 3.4898 | 3.6396 | 3.7890 | Ave | | 3.5024 | | | | 7.8 | | 15.0 | | | |
| 1,3,5-Trichlorobenzene | 1.2561<br>1.2723 | 1.2439 | 1.2909 | 1.2690 | 1.3139 | Ave | | 1.2743 | | | | 2.0 | | 15.0 | | | |
| 2-Methylnaphthalene | 1.1675<br>1.4101 | 1.2466 | 1.4182 | 1.5971 | 1.7053 | Ave | | 1.4241 | | | | 14.3 | | 15.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204067/14 Calibration Date: 10/28/2015 15:21

Instrument ID: A3UX15 Calib Start Date: 10/28/2015 13:05

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/28/2015 14:59

Lab File ID: UXC3830.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.4156 | 0.4225 | 0.1000 | 0.0102 | 0.0100 | 1.7 | 20.0 |
| Chloromethane | Ave | 0.3011 | 0.3110 | 0.1000 | 0.0103 | 0.0100 | 3.3 | 20.0 |
| Vinyl chloride | Ave | 0.3298 | 0.3307 | 0.1000 | 0.0100 | 0.0100 | 0.3 | 20.0 |
| Butadiene | Ave | 0.2931 | 0.2931 | | 0.0100 | 0.0100 | 0.0 | 20.0 |
| Bromomethane | Ave | 0.1790 | 0.1830 | 0.0500 | 0.0102 | 0.0100 | 2.2 | 20.0 |
| Chloroethane | Ave | 0.1902 | 0.1873 | 0.0500 | 0.00985 | 0.0100 | -1.5 | 20.0 |
| Dichlorofluoromethane | Ave | 0.5274 | 0.5397 | | 0.0102 | 0.0100 | 2.3 | 20.0 |
| Trichlorofluoromethane | Ave | 0.4637 | 0.4802 | 0.1000 | 0.0104 | 0.0100 | 3.5 | 20.0 |
| Ethyl ether | Ave | 0.2188 | 0.2222 | | 0.0102 | 0.0100 | 1.6 | 20.0 |
| Acrolein | Ave | 0.0206 | 0.0213 | | 0.0519 | 0.0500 | 3.7 | 20.0 |
| 1,1-Dichloroethene | Ave | 0.3129 | 0.3193 | 0.1000 | 0.0102 | 0.0100 | 2.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2564 | 0.2516 | 0.0500 | 0.00981 | 0.0100 | -1.9 | 20.0 |
| Acetone | Lin1 | | 0.0720 | 0.0500 | 0.0179 | 0.0200 | -10.4 | 20.0 |
| Iodomethane | Ave | 0.4880 | 0.5249 | | 0.0108 | 0.0100 | 7.6 | 20.0 |
| Carbon disulfide | Ave | 0.8576 | 0.9130 | 0.1000 | 0.0106 | 0.0100 | 6.5 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1725 | 0.1700 | | 0.00986 | 0.0100 | -1.4 | 20.0 |
| Methyl acetate | Ave | 0.1509 | 0.1547 | 0.1000 | 0.0513 | 0.0500 | 2.5 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3809 | 0.1000 | 0.0108 | 0.0100 | 8.1 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0133 | 0.0127 | | 0.0952 | 0.100 | -4.8 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7657 | 0.8017 | 0.1000 | 0.0105 | 0.0100 | 4.7 | 20.0 |
| Acrylonitrile | Ave | 0.0817 | 0.0848 | | 0.104 | 0.100 | 3.7 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.3449 | 0.3632 | 0.1000 | 0.0105 | 0.0100 | 5.3 | 20.0 |
| Hexane | Ave | 0.0890 | 0.0954 | | 0.0107 | 0.0100 | 7.2 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.5976 | 0.6462 | 0.1000 | 0.0108 | 0.0100 | 8.1 | 20.0 |
| Vinyl acetate | Ave | 0.3243 | 0.3397 | | 0.0105 | 0.0100 | 4.7 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.3227 | 0.3284 | | 0.0102 | 0.0100 | 1.8 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3710 | 0.3792 | 0.1000 | 0.0102 | 0.0100 | 2.2 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0913 | 0.0904 | 0.0500 | 0.0198 | 0.0200 | -0.9 | 20.0 |
| Bromochloromethane | Ave | 0.1664 | 0.1732 | | 0.0104 | 0.0100 | 4.1 | 20.0 |
| Tetrahydrofuran | Ave | 0.0518 | 0.0521 | | 0.0201 | 0.0200 | 0.6 | 20.0 |
| Chloroform | Ave | 0.6432 | 0.6593 | 0.2000 | 0.0103 | 0.0100 | 2.5 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.4847 | 0.5200 | 0.1000 | 0.0107 | 0.0100 | 7.3 | 20.0 |
| Cyclohexane | Ave | 0.5639 | 0.5858 | 0.1000 | 0.0104 | 0.0100 | 3.9 | 20.0 |
| Carbon tetrachloride | Ave | 0.4107 | 0.4418 | 0.1000 | 0.0108 | 0.0100 | 7.6 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.4747 | 0.5316 | | 0.0112 | 0.0100 | 12.0 | 20.0 |
| Isobutyl alcohol | Ave | 0.0049 | 0.0048 | | 0.240 | 0.250 | -3.8 | 20.0 |
| Benzene | Ave | 1.413 | 1.418 | 0.5000 | 0.0100 | 0.0100 | 0.3 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.5438 | 0.5527 | 0.1000 | 0.0102 | 0.0100 | 1.6 | 20.0 |
| n-Heptane | Ave | 0.0834 | 0.0834 | | 0.0100 | 0.0100 | 0.0 | 20.0 |
| Trichloroethene | Ave | 0.3644 | 0.3846 | 0.2000 | 0.0106 | 0.0100 | 5.5 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204067/14  Calibration Date: 10/28/2015 15:21

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 13:05

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 14:59

Lab File ID: UXC3830.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.5417 | 0.5577 | 0.1000 | 0.0103 | 0.0100 | 2.9 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2981 | 0.3108 | 0.1000 | 0.0104 | 0.0100 | 4.3 | 20.0 |
| Dibromomethane | Ave | 0.1769 | 0.1801 | | 0.0102 | 0.0100 | 1.8 | 20.0 |
| 1,4-Dioxane | Ave | 0.0026 | 0.0022 | | 0.173 | 0.200 | -13.3 | 20.0 |
| Bromodichloromethane | Ave | 0.3932 | 0.3997 | 0.1000 | 0.0102 | 0.0100 | 1.7 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.1542 | 0.1535 | | 0.00995 | 0.0100 | -0.5 | 20.0 |
| cis-1,3-Dichloropropene | Lin1 | | 0.4542 | 0.1500 | 0.00976 | 0.0100 | -2.4 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.1647 | 0.1659 | 0.0500 | 0.0201 | 0.0200 | 0.7 | 20.0 |
| Toluene | Ave | 1.969 | 2.020 | 0.4000 | 0.0103 | 0.0100 | 2.6 | 20.0 |
| trans-1,3-Dichloropropene | Lin1 | | 0.4894 | 0.1000 | 0.00966 | 0.0100 | -3.4 | 20.0 |
| Ethyl methacrylate | Lin1 | | 0.3989 | | 0.00973 | 0.0100 | -2.7 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.3317 | 0.3345 | 0.1000 | 0.0101 | 0.0100 | 0.8 | 20.0 |
| Tetrachloroethene | Ave | 0.3574 | 0.3694 | 0.2000 | 0.0103 | 0.0100 | 3.4 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.6349 | 0.6136 | | 0.00966 | 0.0100 | -3.4 | 20.0 |
| 2-Hexanone | Ave | 0.1536 | 0.1464 | 0.0500 | 0.0191 | 0.0200 | -4.7 | 20.0 |
| Dibromochloromethane | Lin1 | | 0.3187 | | 0.00900 | 0.0100 | -10.0 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.3251 | 0.3340 | | 0.0103 | 0.0100 | 2.7 | 20.0 |
| Chlorobenzene | Ave | 1.266 | 1.274 | 0.3000 | 0.0101 | 0.0100 | 0.6 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3736 | 0.3910 | | 0.0105 | 0.0100 | 4.6 | 20.0 |
| Ethylbenzene | Ave | 0.6530 | 0.6716 | | 0.0103 | 0.0100 | 2.9 | 20.0 |
| m-Xylene & p-Xylene | Ave | 1.789 | 1.810 | | 0.0101 | 0.0100 | 1.2 | 20.0 |
| o-Xylene | Ave | 0.7816 | 0.8009 | | 0.0102 | 0.0100 | 2.5 | 20.0 |
| Styrene | Ave | 1.363 | 1.339 | 0.3000 | 0.00982 | 0.0100 | -1.8 | 20.0 |
| Bromoform | Lin1 | | 0.1536 | 0.1000 | 0.00884 | 0.0100 | -11.6 | 20.0 |
| Isopropylbenzene | Ave | 2.107 | 2.197 | 0.1000 | 0.0104 | 0.0100 | 4.3 | 20.0 |
| Bromobenzene | Ave | 0.9352 | 0.9714 | | 0.0104 | 0.0100 | 3.9 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.6716 | 0.6952 | 0.3000 | 0.0104 | 0.0100 | 3.5 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.2436 | 0.2556 | | 0.0105 | 0.0100 | 4.9 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.1999 | 0.2133 | | 0.0107 | 0.0100 | 6.7 | 20.0 |
| N-Propylbenzene | Ave | 1.009 | 1.078 | | 0.0107 | 0.0100 | 6.8 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8945 | 0.9279 | | 0.0104 | 0.0100 | 3.7 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.344 | 3.567 | | 0.0107 | 0.0100 | 6.7 | 20.0 |
| 4-Chlorotoluene | Ave | 3.463 | 3.593 | | 0.0104 | 0.0100 | 3.7 | 20.0 |
| tert-Butylbenzene | Ave | 2.784 | 2.904 | | 0.0104 | 0.0100 | 4.3 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.455 | 3.643 | | 0.0105 | 0.0100 | 5.4 | 20.0 |
| sec-Butylbenzene | Ave | 3.779 | 3.924 | | 0.0104 | 0.0100 | 3.8 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.848 | 1.859 | 0.6000 | 0.0101 | 0.0100 | 0.6 | 20.0 |
| p-Isopropyltoluene | Ave | 3.438 | 3.575 | | 0.0104 | 0.0100 | 4.0 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.910 | 1.909 | 0.5000 | 0.00999 | 0.0100 | -0.0 | 20.0 |
| n-Butylbenzene | Ave | 2.836 | 2.887 | | 0.0102 | 0.0100 | 1.8 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.764 | 1.756 | 0.4000 | 0.00996 | 0.0100 | -0.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204067/14 Calibration Date: 10/28/2015 15:21

Instrument ID: A3UX15 Calib Start Date: 10/28/2015 13:05

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/28/2015 14:59

Lab File ID: UXC3830.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Lin1 | | 0.0997 | 0.0500 | 0.00916 | 0.0100 | -8.4 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 1.161 | 1.106 | 0.2000 | 0.00952 | 0.0100 | -4.8 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5590 | 0.5093 | | 0.00911 | 0.0100 | -8.9 | 20.0 |
| Naphthalene | Ave | 2.480 | 2.519 | | 0.0102 | 0.0100 | 1.6 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 1.053 | 1.024 | | 0.00972 | 0.0100 | -2.8 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.3239 | 0.3161 | | 0.00715 | 0.00733 | -2.4 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.4605 | 0.4381 | | 0.00697 | 0.00733 | -4.9 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.768 | 1.697 | | 0.00704 | 0.00733 | -4.0 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 0.6700 | 0.6795 | | 0.00744 | 0.00733 | 1.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204067/15  Calibration Date: 10/28/2015 18:00

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 15:44

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 17:37

Lab File ID: UXC3837.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0228 | 0.0226 | | 0.0993 | 0.100 | -0.7 | 20.0 |
| Diisopropyl ether | Ave | 0.2803 | 0.2896 | | 0.0103 | 0.0100 | 3.3 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.5186 | 0.5418 | | 0.0104 | 0.0100 | 4.5 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8119 | 0.8342 | | 0.0103 | 0.0100 | 2.7 | 20.0 |
| Ethyl acetate | Ave | 0.1644 | 0.1655 | | 0.0201 | 0.0200 | 0.7 | 20.0 |
| Propionitrile | Ave | 0.0293 | 0.0301 | | 0.103 | 0.100 | 2.7 | 20.0 |
| Methacrylonitrile | Ave | 0.1324 | 0.1374 | | 0.104 | 0.100 | 3.8 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.6462 | 0.6564 | | 0.0102 | 0.0100 | 1.6 | 20.0 |
| n-Butanol | Ave | 0.0035 | 0.0031 | | 0.219 | 0.250 | -12.3 | 20.0 |
| Methyl methacrylate | Ave | 0.1803 | 0.1875 | | 0.0208 | 0.0200 | 4.0 | 20.0 |
| 2-Nitropropane | Ave | 0.0274 | 0.0237 | | 0.0173 | 0.0200 | -13.7 | 20.0 |
| 1-Chlorohexane | Ave | 0.5138 | 0.5158 | | 0.0100 | 0.0100 | 0.4 | 20.0 |
| Cyclohexanone | Ave | 0.0128 | 0.0120 | | 0.0933 | 0.100 | -6.7 | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.502 | 3.574 | | 0.0102 | 0.0100 | 2.0 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.274 | 1.242 | | 0.00974 | 0.0100 | -2.6 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.424 | 1.167 | | 0.0164 | 0.0200 | -18.1 | 20.0 |

FORM V
GC/MS VOA INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: BFB472.D BFB Injection Date: 11/12/2015

Instrument ID: A3UX15 BFB Injection Time: 10:14

Analysis Batch No.: 206461

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 50 | 15.0 - 40.0 % of mass 95 | 16.9 | |
| 75 | 30.0 - 60.0 % of mass 95 | 53.2 | |
| 95 | Base Peak, 100% relative abundance | 100.0 | |
| 96 | 5.0 - 9.0 % of mass 95 | 6.7 | |
| 173 | Less than 2.0 % of mass 174 | 0.6 | (0.8)1 |
| 174 | 50.0 - 120.00 % of mass 95 | 80.0 | |
| 175 | 5.0 - 9.0 % of mass 174 | 5.8 | (7.3)1 |
| 176 | 95.0 - 101.0 % of mass 174 | 78.1 | (97.7)1 |
| 177 | 5.0 - 9.0 % of mass 176 | 5.2 | (6.7)2 |

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCVIS 240-206461/2 | UXC4392.D | 11/12/2015 | 10:41 |
| | LCS 240-206461/4 | UXC4393.D | 11/12/2015 | 11:03 |
| | CCV 240-206461/3 | UXC4394.D | 11/12/2015 | 11:26 |
| | MB 240-206461/6 | UXC4396.D | 11/12/2015 | 12:11 |
| TB-110515 | 240-57597-16 | UXC4397.D | 11/12/2015 | 12:33 |
| | 240-57564-B-1 MS | UXC4417.D | 11/12/2015 | 20:05 |
| | 240-57564-B-1 MSD | UXC4418.D | 11/12/2015 | 20:27 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206461/2 Calibration Date: 11/12/2015 10:41

Instrument ID: A3UX15 Calib Start Date: 10/28/2015 13:05

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/28/2015 14:59

Lab File ID: UXC4392.D Conc. Units: ng/uL Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Dichlorodifluoromethane | Ave | 0.4156 | 0.4748 | 0.1000 | 0.0114 | 0.0100 | 14.3 | 20.0 |
| Chloromethane | Ave | 0.3011 | 0.2531 | 0.1000 | 0.00841 | 0.0100 | -15.9 | 20.0 |
| Vinyl chloride | Ave | 0.3298 | 0.2842 | 0.1000 | 0.00862 | 0.0100 | -13.8 | 20.0 |
| Butadiene | Ave | 0.2931 | 0.2523 | | 0.00861 | 0.0100 | -13.9 | 20.0 |
| Bromomethane | Ave | 0.1790 | 0.1968 | 0.0500 | 0.0110 | 0.0100 | 10.0 | 20.0 |
| Chloroethane | Ave | 0.1902 | 0.1836 | 0.0500 | 0.00965 | 0.0100 | -3.5 | 20.0 |
| Dichlorofluoromethane | Ave | 0.5274 | 0.5573 | | 0.0106 | 0.0100 | 5.7 | 20.0 |
| Trichlorofluoromethane | Ave | 0.4637 | 0.4470 | 0.1000 | 0.00964 | 0.0100 | -3.6 | 20.0 |
| Ethyl ether | Ave | 0.2188 | 0.1979 | | 0.00905 | 0.0100 | -9.5 | 20.0 |
| Acrolein | Ave | 0.0206 | 0.0157 | | 0.0381 | 0.0500 | -23.9* | 20.0 |
| 1,1-Dichloroethene | Ave | 0.3129 | 0.3099 | 0.1000 | 0.00990 | 0.0100 | -1.0 | 20.0 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane | Ave | 0.2564 | 0.2706 | 0.0500 | 0.0106 | 0.0100 | 5.5 | 20.0 |
| Acetone | Lin1 | | 0.0764 | 0.0500 | 0.0191 | 0.0200 | -4.3 | 20.0 |
| Iodomethane | Ave | 0.4880 | 0.5373 | | 0.0110 | 0.0100 | 10.1 | 20.0 |
| Carbon disulfide | Ave | 0.8576 | 0.8155 | 0.1000 | 0.00951 | 0.0100 | -4.9 | 20.0 |
| 3-Chloro-1-propene | Ave | 0.1725 | 0.1609 | | 0.00933 | 0.0100 | -6.7 | 20.0 |
| Methyl acetate | Ave | 0.1509 | 0.1291 | 0.1000 | 0.0428 | 0.0500 | -14.4 | 20.0 |
| Methylene Chloride | Lin1 | | 0.3455 | 0.1000 | 0.00970 | 0.0100 | -3.0 | 20.0 |
| tert-Butyl alcohol (TBA) | Ave | 0.0133 | 0.0120 | | 0.0900 | 0.100 | -10.0 | 20.0 |
| Methyl tert-butyl ether | Ave | 0.7657 | 0.7865 | 0.1000 | 0.0103 | 0.0100 | 2.7 | 20.0 |
| trans-1,2-Dichloroethene | Ave | 0.3449 | 0.3440 | 0.1000 | 0.00998 | 0.0100 | -0.2 | 20.0 |
| Acrylonitrile | Ave | 0.0817 | 0.0722 | | 0.0883 | 0.100 | -11.7 | 20.0 |
| Hexane | Ave | 0.0890 | 0.0821 | | 0.00922 | 0.0100 | -7.8 | 20.0 |
| 1,1-Dichloroethane | Ave | 0.5976 | 0.5650 | 0.1000 | 0.00945 | 0.0100 | -5.5 | 20.0 |
| Vinyl acetate | Ave | 0.3243 | 0.3339 | | 0.0103 | 0.0100 | 2.9 | 20.0 |
| 2,2-Dichloropropane | Ave | 0.3227 | 0.3420 | | 0.0106 | 0.0100 | 6.0 | 20.0 |
| cis-1,2-Dichloroethene | Ave | 0.3710 | 0.3555 | 0.1000 | 0.00958 | 0.0100 | -4.2 | 20.0 |
| 2-Butanone (MEK) | Ave | 0.0913 | 0.0782 | 0.0500 | 0.0171 | 0.0200 | -14.3 | 20.0 |
| Bromochloromethane | Ave | 0.1664 | 0.1838 | | 0.0110 | 0.0100 | 10.5 | 20.0 |
| Tetrahydrofuran | Ave | 0.0518 | 0.0467 | | 0.0181 | 0.0200 | -9.7 | 20.0 |
| Chloroform | Ave | 0.6432 | 0.6727 | 0.2000 | 0.0105 | 0.0100 | 4.6 | 20.0 |
| 1,1,1-Trichloroethane | Ave | 0.4847 | 0.5235 | 0.1000 | 0.0108 | 0.0100 | 8.0 | 20.0 |
| Cyclohexane | Ave | 0.5639 | 0.4901 | 0.1000 | 0.00869 | 0.0100 | -13.1 | 20.0 |
| 1,1-Dichloropropene | Ave | 0.4747 | 0.4704 | | 0.00991 | 0.0100 | -0.9 | 20.0 |
| Carbon tetrachloride | Ave | 0.4107 | 0.4780 | 0.1000 | 0.0116 | 0.0100 | 16.4 | 20.0 |
| Isobutyl alcohol | Ave | 0.0049 | 0.0045 | | 0.228 | 0.250 | -8.9 | 20.0 |
| Benzene | Ave | 1.413 | 1.334 | 0.5000 | 0.00944 | 0.0100 | -5.6 | 20.0 |
| 1,2-Dichloroethane | Ave | 0.5438 | 0.5731 | 0.1000 | 0.0105 | 0.0100 | 5.4 | 20.0 |
| n-Heptane | Ave | 0.0834 | 0.0828 | | 0.00993 | 0.0100 | -0.7 | 20.0 |
| Trichloroethene | Ave | 0.3644 | 0.3950 | 0.2000 | 0.0108 | 0.0100 | 8.4 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206461/2  Calibration Date: 11/12/2015 10:41

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 13:05

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 14:59

Lab File ID: UXC4392.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Methylcyclohexane | Ave | 0.5417 | 0.4958 | 0.1000 | 0.00915 | 0.0100 | -8.5 | 20.0 |
| 1,2-Dichloropropane | Ave | 0.2981 | 0.2720 | 0.1000 | 0.00913 | 0.0100 | -8.7 | 20.0 |
| Dibromomethane | Ave | 0.1769 | 0.1852 | | 0.0105 | 0.0100 | 4.7 | 20.0 |
| 1,4-Dioxane | Ave | 0.0026 | 0.0024 | | 0.187 | 0.200 | -6.5 | 20.0 |
| Bromodichloromethane | Ave | 0.3932 | 0.4333 | 0.1000 | 0.0110 | 0.0100 | 10.2 | 20.0 |
| 2-Chloroethyl vinyl ether | Ave | 0.1542 | 0.1349 | | 0.0175 | 0.0200 | -12.5 | 20.0 |
| cis-1,3-Dichloropropene | Lin1 | | 0.4161 | 0.1500 | 0.00898 | 0.0100 | -10.2 | 20.0 |
| 4-Methyl-2-pentanone (MIBK) | Ave | 0.1647 | 0.1558 | 0.0500 | 0.0189 | 0.0200 | -5.4 | 20.0 |
| Toluene | Ave | 1.969 | 1.768 | 0.4000 | 0.00898 | 0.0100 | -10.2 | 20.0 |
| trans-1,3-Dichloropropene | Lin1 | | 0.4223 | 0.1000 | 0.00842 | 0.0100 | -15.8 | 20.0 |
| Ethyl methacrylate | Lin1 | | 0.3389 | | 0.00834 | 0.0100 | -16.6 | 20.0 |
| 1,1,2-Trichloroethane | Ave | 0.3317 | 0.3105 | 0.1000 | 0.00936 | 0.0100 | -6.4 | 20.0 |
| Tetrachloroethene | Ave | 0.3574 | 0.3563 | 0.2000 | 0.00997 | 0.0100 | -0.3 | 20.0 |
| 1,3-Dichloropropane | Ave | 0.6349 | 0.5496 | | 0.00866 | 0.0100 | -13.4 | 20.0 |
| 2-Hexanone | Ave | 0.1536 | 0.1235 | 0.0500 | 0.0161 | 0.0200 | -19.6 | 20.0 |
| Dibromochloromethane | Lin1 | | 0.3343 | | 0.00941 | 0.0100 | -5.9 | 20.0 |
| 1,2-Dibromoethane | Ave | 0.3251 | 0.3105 | | 0.00955 | 0.0100 | -4.5 | 20.0 |
| Chlorobenzene | Ave | 1.266 | 1.196 | 0.3000 | 0.00944 | 0.0100 | -5.6 | 20.0 |
| 1,1,1,2-Tetrachloroethane | Ave | 0.3736 | 0.4061 | | 0.0109 | 0.0100 | 8.7 | 20.0 |
| Ethylbenzene | Ave | 0.6530 | 0.6239 | | 0.00956 | 0.0100 | -4.4 | 20.0 |
| m-Xylene & p-Xylene | Ave | 1.789 | 1.688 | | 0.00943 | 0.0100 | -5.7 | 20.0 |
| o-Xylene | Ave | 0.7816 | 0.7445 | | 0.00952 | 0.0100 | -4.8 | 20.0 |
| Styrene | Ave | 1.363 | 1.327 | 0.3000 | 0.00974 | 0.0100 | -2.6 | 20.0 |
| Bromoform | Lin1 | | 0.1723 | 0.1000 | 0.00983 | 0.0100 | -1.7 | 20.0 |
| Isopropylbenzene | Ave | 2.107 | 2.063 | 0.1000 | 0.00979 | 0.0100 | -2.1 | 20.0 |
| Bromobenzene | Ave | 0.9352 | 0.8652 | | 0.00925 | 0.0100 | -7.5 | 20.0 |
| 1,1,2,2-Tetrachloroethane | Ave | 0.6716 | 0.5654 | 0.3000 | 0.00842 | 0.0100 | -15.8 | 20.0 |
| 1,2,3-Trichloropropane | Ave | 0.2436 | 0.2253 | | 0.00925 | 0.0100 | -7.5 | 20.0 |
| trans-1,4-Dichloro-2-butene | Ave | 0.1999 | 0.1452 | | 0.00726 | 0.0100 | -27.4* | 20.0 |
| N-Propylbenzene | Ave | 1.009 | 0.8864 | | 0.00878 | 0.0100 | -12.2 | 20.0 |
| 2-Chlorotoluene | Ave | 0.8945 | 0.7988 | | 0.00893 | 0.0100 | -10.7 | 20.0 |
| 1,3,5-Trimethylbenzene | Ave | 3.344 | 2.940 | | 0.00879 | 0.0100 | -12.1 | 20.0 |
| 4-Chlorotoluene | Ave | 3.463 | 2.986 | | 0.00862 | 0.0100 | -13.8 | 20.0 |
| tert-Butylbenzene | Ave | 2.784 | 2.450 | | 0.00880 | 0.0100 | -12.0 | 20.0 |
| 1,2,4-Trimethylbenzene | Ave | 3.455 | 3.098 | | 0.00897 | 0.0100 | -10.3 | 20.0 |
| sec-Butylbenzene | Ave | 3.779 | 3.234 | | 0.00856 | 0.0100 | -14.4 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.848 | 1.686 | 0.6000 | 0.00912 | 0.0100 | -8.8 | 20.0 |
| p-Isopropyltoluene | Ave | 3.438 | 3.118 | | 0.00907 | 0.0100 | -9.3 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.910 | 1.691 | 0.5000 | 0.00885 | 0.0100 | -11.5 | 20.0 |
| n-Butylbenzene | Ave | 2.836 | 2.292 | | 0.00808 | 0.0100 | -19.2 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.764 | 1.567 | 0.4000 | 0.00889 | 0.0100 | -11.1 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206461/2  Calibration Date: 11/12/2015 10:41

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 13:05

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 14:59

Lab File ID: UXC4392.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,2-Dibromo-3-Chloropropane | Lin1 | | 0.0997 | 0.0500 | 0.00917 | 0.0100 | -8.3 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 1.161 | 1.004 | 0.2000 | 0.00864 | 0.0100 | -13.6 | 20.0 |
| Hexachlorobutadiene | Ave | 0.5590 | 0.4334 | | 0.00775 | 0.0100 | -22.5* | 20.0 |
| Naphthalene | Ave | 2.480 | 2.098 | | 0.00846 | 0.0100 | -15.4 | 20.0 |
| 1,2,3-Trichlorobenzene | Ave | 1.053 | 0.8905 | | 0.00846 | 0.0100 | -15.4 | 20.0 |
| Dibromofluoromethane (Surr) | Ave | 0.3239 | 0.3247 | | 0.00735 | 0.00733 | 0.2 | 20.0 |
| 1,2-Dichloroethane-d4 (Surr) | Ave | 0.4605 | 0.4500 | | 0.00716 | 0.00733 | -2.3 | 20.0 |
| Toluene-d8 (Surr) | Ave | 1.768 | 1.546 | | 0.00641 | 0.00733 | -12.5 | 20.0 |
| 4-Bromofluorobenzene (Surr) | Ave | 0.6700 | 0.6842 | | 0.00749 | 0.00733 | 2.1 | 20.0 |

FORM VII
GC/MS VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206461/3  Calibration Date: 11/12/2015 11:26

Instrument ID: A3UX15  Calib Start Date: 10/28/2015 15:44

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/28/2015 17:37

Lab File ID: UXC4394.D  Conc. Units: ng/uL  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Acetonitrile | Ave | 0.0228 | 0.0075 | | 0.0330 | 0.100 | -67.0* | 20.0 |
| Diisopropyl ether | Ave | 0.2803 | 0.2793 | | 0.00996 | 0.0100 | -0.4 | 20.0 |
| 2-Chloro-1,3-butadiene | Ave | 0.5186 | 0.4985 | | 0.00961 | 0.0100 | -3.9 | 20.0 |
| Ethyl t-butyl ether | Ave | 0.8119 | 0.7747 | | 0.00954 | 0.0100 | -4.6 | 20.0 |
| Ethyl acetate | Ave | 0.1644 | 0.1216 | | 0.0148 | 0.0200 | -26.0* | 20.0 |
| Propionitrile | Ave | 0.0293 | 0.0194 | | 0.0663 | 0.100 | -33.7* | 20.0 |
| Methacrylonitrile | Ave | 0.1324 | 0.1064 | | 0.0804 | 0.100 | -19.6 | 20.0 |
| Tert-amyl methyl ether | Ave | 0.6462 | 0.6591 | | 0.0102 | 0.0100 | 2.0 | 20.0 |
| n-Butanol | Ave | 0.0035 | 0.0016 | | 0.115 | 0.250 | -54.1* | 20.0 |
| Methyl methacrylate | Ave | 0.1803 | 0.1522 | | 0.0169 | 0.0200 | -15.6 | 20.0 |
| 2-Nitropropane | Ave | 0.0274 | 0.0252 | | 0.0184 | 0.0200 | -8.2 | 20.0 |
| 1-Chlorohexane | Ave | 0.5138 | 0.4605 | | 0.00896 | 0.0100 | -10.4 | 20.0 |
| Cyclohexanone | Ave | 0.0128 | 0.0048 | | 0.0378 | 0.100 | -62.2* | 20.0 |
| 1,2,3-Trimethylbenzene | Ave | 3.502 | 3.168 | | 0.00905 | 0.0100 | -9.5 | 20.0 |
| 1,3,5-Trichlorobenzene | Ave | 1.274 | 1.080 | | 0.00847 | 0.0100 | -15.3 | 20.0 |
| 2-Methylnaphthalene | Ave | 1.424 | 0.9190 | | 0.0129 | 0.0200 | -35.5* | 20.0 |

FORM II
GC/MS SEMI VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low

GC Column (1): DB-624  ID: 0.18(mm)

| Client Sample ID | Lab Sample ID | 2FP # | PHL # | NBZ # | FBP # | TBP # | TPH # |
|---|---|---|---|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 64 | 65 | 62 | 67 | 44 | 76 |
| F-VB-UST-2-14 | 240-57597-2 | 61 | 61 | 58 | 63 | 42 | 78 |
| F-VB-UST-3-14 | 240-57597-3 | 61 | 60 | 58 | 63 | 41 | 73 |
| F-VB-UST-4-14 | 240-57597-4 | 72 | 72 | 69 | 72 | 42 | 85 |
| F-VB-UST-5-14 | 240-57597-5 | 68 | 67 | 64 | 69 | 25 | 77 |
| F-VB-UST-6-14 | 240-57597-6 | 62 | 60 | 62 | 68 | 33 | 79 |
| F-VB-UST-7-14 | 240-57597-7 | 67 | 64 | 62 | 70 | 38 | 80 |
| F-VB-UST-8-14 | 240-57597-8 | 56 | 54 | 52 | 56 | 32 | 70 |
| F-VB-UST-9-14 | 240-57597-9 | 71 | 68 | 68 | 73 | 45 | 87 |
| F-VB-UST-10-14 | 240-57597-10 | 59 | 61 | 60 | 64 | 40 | 78 |
| F-VB-UST-11-14 | 240-57597-11 | 78 | 80 | 80 | 73 | 40 | 88 |
| F-VB-UST-12-14 | 240-57597-12 | 67 | 64 | 67 | 63 | 32 | 72 |
| F-VB-UST-13-14 | 240-57597-13 | 65 | 57 | 65 | 59 | 33 | 77 |
| F-VB-UST-14-14 | 240-57597-14 | 70 | 62 | 73 | 68 | 28 | 79 |
| F-VB-UST-DUP | 240-57597-15 | 71 | 60 | 74 | 68 | 24 | 81 |
| | MB 240-205786/23-A | 66 | 66 | 65 | 68 | 48 | 80 |
| | MB 240-205866/8-A | 81 | 88 | 78 | 76 | 49 | 87 |
| | LCS 240-205786/24-A | 79 | 79 | 79 | 81 | 44 | 93 |
| | LCS 240-205866/9-A | 72 | 75 | 76 | 64 | 33 | 76 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 69 | 62 | 78 | 71 | 30 | 80 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 61 | 53 | 65 | 60 | 36 | 77 |

| | QC LIMITS |
|---|---|
| 2FP = 2-Fluorophenol (Surr) | 24-110 |
| PHL = Phenol-d5 (Surr) | 26-110 |
| NBZ = Nitrobenzene-d5 (Surr) | 20-110 |
| FBP = 2-Fluorobiphenyl (Surr) | 24-110 |
| TBP = 2,4,6-Tribromophenol (Surr) | 10-110 |
| TPH = Terphenyl-d14 (Surr) | 36-110 |

# Column to be used to flag recovery values

FORM II 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: STD6 240-200091/10  Date Analyzed: 10/02/2015 14:04

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 51002010.D  Heated Purge: (Y/N) N

Calibration ID: 30980

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 221463 | 6.23 | 906472 | 7.34 | 529260 | 8.85 |
| UPPER LIMIT | | 442926 | 6.73 | 1812944 | 7.84 | 1058520 | 9.35 |
| LOWER LIMIT | | 110732 | 5.73 | 453236 | 6.84 | 264630 | 8.35 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-200091/11 | | 242280 | 6.23 | 1000939 | 7.34 | 580042 | 8.85 |
| CCV 240-206256/2 CCVIS | | 183138 | 6.15 | 734641 | 7.27 | 425153 | 8.77 |

DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Sample No.: STD6 240-200091/10 Date Analyzed: 10/02/2015 14:04

Instrument ID: A4HP10 GC Column: DB-624 ID: 0.18(mm)

Lab File ID (Standard): 51002010.D Heated Purge: (Y/N) N

Calibration ID: 30980

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 840219 | 10.12 | 1037091 | 12.82 | 955028 | 15.02 |
| UPPER LIMIT | | 1680438 | 10.62 | 2074182 | 13.32 | 1910056 | 15.52 |
| LOWER LIMIT | | 420110 | 9.62 | 518546 | 12.32 | 477514 | 14.52 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-200091/11 | | 929663 | 10.12 | 1154414 | 12.82 | 1074213 | 15.02 |
| CCV 240-206256/2 CCVIS | | 678359 | 10.04 | 794127 | 12.71 | 678609 | 14.89 |

PHN = Phenanthrene-d10
CRY = Chrysene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206256/2  Date Analyzed: 11/11/2015 09:34

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 51111002.D  Heated Purge: (Y/N) N

Calibration ID: 31049

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 183138 | 6.15 | 734641 | 7.27 | 425153 | 8.77 |
| UPPER LIMIT | | 366276 | 6.65 | 1469282 | 7.77 | 850306 | 9.27 |
| LOWER LIMIT | | 91569 | 5.65 | 367321 | 6.77 | 212577 | 8.27 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-205786/23-A | | 169364 | 6.16 | 695817 | 7.27 | 416939 | 8.77 |
| LCS 240-205786/24-A | | 167619 | 6.16 | 683191 | 7.27 | 403392 | 8.77 |
| 240-57597-1 | F-VB-UST-1-14 | 170794 | 6.16 | 710718 | 7.27 | 416040 | 8.77 |
| 240-57597-2 | F-VB-UST-2-14 | 172301 | 6.16 | 704300 | 7.27 | 408420 | 8.77 |
| 240-57597-3 | F-VB-UST-3-14 | 168152 | 6.16 | 687534 | 7.27 | 389818 | 8.77 |
| 240-57597-4 | F-VB-UST-4-14 | 159703 | 6.16 | 681534 | 7.27 | 397658 | 8.77 |
| 240-57597-5 | F-VB-UST-5-14 | 169966 | 6.16 | 709398 | 7.27 | 411661 | 8.77 |
| 240-57597-6 | F-VB-UST-6-14 | 173464 | 6.16 | 694821 | 7.27 | 402882 | 8.77 |
| 240-57597-7 | F-VB-UST-7-14 | 166312 | 6.16 | 698152 | 7.27 | 395626 | 8.77 |
| 240-57597-8 | F-VB-UST-8-14 | 170392 | 6.16 | 690624 | 7.27 | 399583 | 8.77 |
| 240-57597-9 | F-VB-UST-9-14 | 167764 | 6.16 | 674344 | 7.27 | 396455 | 8.77 |
| 240-57597-10 | F-VB-UST-10-14 | 165361 | 6.16 | 676786 | 7.27 | 404905 | 8.77 |

DCB = 1,4-Dichlorobenzene-d4
DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
NPT = Naphthalene-d8
Area Limit = 50%-200% of internal standard area
ANT = Acenaphthene-d10
RT Limit = ± 0.5 minutes of internal standard RT
ANT = Acenaphthene-d10

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206256/2  Date Analyzed: 11/11/2015 09:34

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)

Lab File ID (Standard): 51111002.D  Heated Purge: (Y/N) N

Calibration ID: 31049

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 678359 | 10.04 | 794127 | 12.71 | 678609 | 14.89 |
| UPPER LIMIT | | 1356718 | 10.54 | 1588254 | 13.21 | 1357218 | 15.39 |
| LOWER LIMIT | | 339180 | 9.54 | 397064 | 12.21 | 339305 | 14.39 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-205786/23-A | | 633596 | 10.04 | 730132 | 12.70 | 644324 | 14.88 |
| LCS 240-205786/24-A | | 651601 | 10.04 | 738522 | 12.70 | 635222 | 14.89 |
| 240-57597-1 | F-VB-UST-1-14 | 661549 | 10.04 | 736014 | 12.70 | 508559 | 14.88 |
| 240-57597-2 | F-VB-UST-2-14 | 651468 | 10.04 | 719477 | 12.70 | 573126 | 14.88 |
| 240-57597-3 | F-VB-UST-3-14 | 616636 | 10.04 | 695392 | 12.70 | 523351 | 14.88 |
| 240-57597-4 | F-VB-UST-4-14 | 624736 | 10.04 | 703819 | 12.70 | 478985 | 14.88 |
| 240-57597-5 | F-VB-UST-5-14 | 639217 | 10.04 | 741747 | 12.70 | 497306 | 14.88 |
| 240-57597-6 | F-VB-UST-6-14 | 640684 | 10.04 | 728874 | 12.70 | 488005 | 14.88 |
| 240-57597-7 | F-VB-UST-7-14 | 639827 | 10.04 | 723063 | 12.70 | 509629 | 14.88 |
| 240-57597-8 | F-VB-UST-8-14 | 631926 | 10.04 | 719274 | 12.70 | 575492 | 14.88 |
| 240-57597-9 | F-VB-UST-9-14 | 624386 | 10.04 | 704835 | 12.70 | 495093 | 14.88 |
| 240-57597-10 | F-VB-UST-10-14 | 640666 | 10.04 | 718652 | 12.70 | 538508 | 14.88 |

PHN = Phenanthrene-d10
PHN = Phenanthrene-d10
CRY = Chrysene-d12
CRY = Chrysene-d12
PRY = Perylene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: STD6 240-204569/10  Date Analyzed: 10/30/2015 16:42

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)

Lab File ID (Standard): 51030010.D  Heated Purge: (Y/N) N

Calibration ID: 31170

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 251143 | 6.42 | 1074814 | 7.54 | 555801 | 9.04 |
| UPPER LIMIT | | 502286 | 6.92 | 2149628 | 8.04 | 1111602 | 9.54 |
| LOWER LIMIT | | 125572 | 5.92 | 537407 | 7.04 | 277901 | 8.54 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-204569/11 | | 291290 | 6.42 | 1233894 | 7.54 | 643573 | 9.04 |
| CCV 240-206233/2 CCVIS | | 263455 | 6.30 | 1198567 | 7.42 | 612244 | 8.92 |

DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Sample No.: STD6 240-204569/10 Date Analyzed: 10/30/2015 16:42

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm)

Lab File ID (Standard): 51030010.D Heated Purge: (Y/N) N

Calibration ID: 31170

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 848550 | 10.31 | 938909 | 12.62 | 798084 | 14.43 |
| UPPER LIMIT | | 1697100 | 10.81 | 1877818 | 13.12 | 1596168 | 14.93 |
| LOWER LIMIT | | 424275 | 9.81 | 469455 | 12.12 | 399042 | 13.93 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-204569/11 | | 1041455 | 10.31 | 1160345 | 12.61 | 991819 | 14.43 |
| CCV 240-206233/2 CCVIS | | 908264 | 10.19 | 962378 | 12.46 | 831716 | 14.18 |

PHN = Phenanthrene-d10
CRY = Chrysene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206233/2  Date Analyzed: 11/11/2015 07:49

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)

Lab File ID (Standard): 51111002.D  Heated Purge: (Y/N) N

Calibration ID: 31170

| | | DCB | | NPT | | ANT | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 263455 | 6.30 | 1198567 | 7.42 | 612244 | 8.92 |
| UPPER LIMIT | | 526910 | 6.80 | 2397134 | 7.92 | 1224488 | 9.42 |
| LOWER LIMIT | | 131728 | 5.80 | 599284 | 6.92 | 306122 | 8.42 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-205866/8-A | | 217654 | 6.30 | 987205 | 7.42 | 480730 | 8.92 |
| LCS 240-205866/9-A | | 227686 | 6.30 | 1006110 | 7.42 | 522813 | 8.92 |
| 240-57597-11 | F-VB-UST-11-14 | 214487 | 6.29 | 930769 | 7.42 | 469347 | 8.92 |
| 240-57597-12 | F-VB-UST-12-14 | 207603 | 6.30 | 930370 | 7.42 | 469544 | 8.92 |
| 240-57597-13 | F-VB-UST-13-14 | 229993 | 6.30 | 1005170 | 7.42 | 509477 | 8.92 |
| 240-57597-14 | F-VB-UST-14-14 | 236787 | 6.30 | 998849 | 7.42 | 507858 | 8.92 |
| 240-57597-15 | F-VB-UST-DUP | 220994 | 6.30 | 932862 | 7.42 | 478231 | 8.92 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 216080 | 6.30 | 960244 | 7.42 | 481718 | 8.92 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 234420 | 6.30 | 1033463 | 7.42 | 522594 | 8.92 |

DCB = 1,4-Dichlorobenzene-d4
DCB = 1,4-Dichlorobenzene-d4
NPT = Naphthalene-d8
NPT = Naphthalene-d8
Area Limit = 50%-200% of internal standard area
ANT = Acenaphthene-d10
RT Limit = ± 0.5 minutes of internal standard RT
ANT = Acenaphthene-d10

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206233/2  Date Analyzed: 11/11/2015 07:49

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)

Lab File ID (Standard): 51111002.D  Heated Purge: (Y/N) N

Calibration ID: 31170

| | | PHN | | CRY | | PRY | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 908264 | 10.19 | 962378 | 12.46 | 831716 | 14.18 |
| UPPER LIMIT | | 1816528 | 10.69 | 1924756 | 12.96 | 1663432 | 14.68 |
| LOWER LIMIT | | 454132 | 9.69 | 481189 | 11.96 | 415858 | 13.68 |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| MB 240-205866/8-A | | 733495 | 10.18 | 742087 | 12.46 | 654968 | 14.18 |
| LCS 240-205866/9-A | | 774857 | 10.19 | 814640 | 12.46 | 725274 | 14.18 |
| 240-57597-11 | F-VB-UST-11-14 | 720283 | 10.19 | 731765 | 12.46 | 528934 | 14.18 |
| 240-57597-12 | F-VB-UST-12-14 | 701609 | 10.19 | 717786 | 12.46 | 585455 | 14.18 |
| 240-57597-13 | F-VB-UST-13-14 | 754672 | 10.19 | 785246 | 12.46 | 684137 | 14.18 |
| 240-57597-14 | F-VB-UST-14-14 | 760631 | 10.19 | 787874 | 12.46 | 628812 | 14.18 |
| 240-57597-15 | F-VB-UST-DUP | 704693 | 10.18 | 734178 | 12.46 | 582028 | 14.17 |
| 240-57597-15 MS | F-VB-UST-DUP MS | 739261 | 10.19 | 786491 | 12.46 | 699060 | 14.18 |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 822314 | 10.19 | 841298 | 12.46 | 788802 | 14.18 |

PHN = Phenanthrene-d10
PHN = Phenanthrene-d10
CRY = Chrysene-d12
CRY = Chrysene-d12
PRY = Perylene-d12
PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8270D

FORM IV
GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab File ID: 51111004.D  Lab Sample ID: MB 240-205786/23-A

Matrix: Solid  Date Extracted: 11/09/2015 06:12

Instrument ID: A4HP10  Date Analyzed: 11/11/2015 10:27

Level:(Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-205786/24-A | 51111005.D | 11/11/2015 10:51 |
| F-VB-UST-1-14 | 240-57597-1 | 51111006.D | 11/11/2015 11:16 |
| F-VB-UST-2-14 | 240-57597-2 | 51111007.D | 11/11/2015 11:40 |
| F-VB-UST-3-14 | 240-57597-3 | 51111008.D | 11/11/2015 12:05 |
| F-VB-UST-4-14 | 240-57597-4 | 51111009.D | 11/11/2015 12:30 |
| F-VB-UST-5-14 | 240-57597-5 | 51111010.D | 11/11/2015 12:54 |
| F-VB-UST-6-14 | 240-57597-6 | 51111011.D | 11/11/2015 13:19 |
| F-VB-UST-7-14 | 240-57597-7 | 51111012.D | 11/11/2015 13:44 |
| F-VB-UST-8-14 | 240-57597-8 | 51111013.D | 11/11/2015 14:08 |
| F-VB-UST-9-14 | 240-57597-9 | 51111014.D | 11/11/2015 14:33 |
| F-VB-UST-10-14 | 240-57597-10 | 51111015.D | 11/11/2015 14:57 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205786/23-A

Matrix: Solid | Lab File ID: 51111004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:27

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 3.5 | U | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 9.5 | U | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 25 | U | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 8.9 | U | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 20 | U | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 20 | U | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 21 | U | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 17 | U | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 21 | U | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 0.45 | U | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 8.2 | U | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 0.50 | U | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 11 | U | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 9.1 | U | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 8.3 | U | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 18 | U | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 16 | U | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 9.2 | U | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 13 | U | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 21 | U | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 17 | U | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 13 | U | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 26 | U | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 17 | U | 330 | 17 |
| 83-32-9 | Acenaphthene | 0.76 | U | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 0.35 | U | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 9.2 | U | 100 | 9.2 |
| 120-12-7 | Anthracene | 0.78 | U | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 9.1 | U | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 12 | U | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 0.63 | U | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 0.64 | U | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 0.59 | U | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.35 | U | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205786/23-A

Matrix: Solid | Lab File ID: 51111004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:27

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.68 | U | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 22 | U | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.0 | U | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 19 | U | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 10 | U | 70 | 10 |
| 105-60-2 | Caprolactam | 37 | U | 330 | 37 |
| 86-74-8 | Carbazole | 27 | U | 50 | 27 |
| 218-01-9 | Chrysene | 1.1 | U | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.66 | U | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 0.66 | U | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 16 | U | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 17 | U | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 15 | U | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 7.9 | U | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 0.55 | U | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 0.53 | U | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 2.1 | U | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 5.6 | U | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 8.1 | U | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 9.0 | U | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.35 | U | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 13 | U | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 6.3 | U | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 21 | U | 50 | 21 |
| 91-20-3 | Naphthalene | 0.82 | U | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 2.2 | U | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 9.1 | U | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 0.73 | U | 6.7 | 0.73 |
| 108-95-2 | Phenol | 7.3 | U | 50 | 7.3 |
| 129-00-0 | Pyrene | 0.44 | U | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 20 | U | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205786/23-A

Matrix: Solid | Lab File ID: 51111004.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 06:12

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 10:27

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206256 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 80 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 66 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 65 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 66 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 68 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 48 | | 10-110 |

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111005.D

Lab ID: LCS 240-205786/24-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 525 | 79 | 35-110 | |
| bis (2-chloroisopropyl) ether | 667 | 611 | 92 | 29-110 | |
| 2,4,5-Trichlorophenol | 667 | 428 | 64 | 25-110 | |
| 2,4,6-Trichlorophenol | 667 | 319 | 48 | 12-110 | |
| 2,4-Dichlorophenol | 667 | 532 | 80 | 39-110 | |
| 2,4-Dimethylphenol | 667 | 462 | 69 | 29-110 | |
| 2,4-Dinitrophenol | 1330 | 481 | 36 | 10-110 | |
| 2,4-Dinitrotoluene | 667 | 605 | 91 | 48-110 | |
| 2,6-Dinitrotoluene | 667 | 563 | 84 | 45-110 | |
| 2-Chloronaphthalene | 667 | 523 | 78 | 32-110 | |
| 2-Chlorophenol | 667 | 528 | 79 | 37-110 | |
| 2-Methylnaphthalene | 667 | 550 | 83 | 36-110 | |
| 2-Methylphenol | 667 | 576 | 86 | 41-110 | |
| 2-Nitroaniline | 667 | 596 | 89 | 45-110 | |
| 2-Nitrophenol | 667 | 584 | 88 | 34-110 | |
| 3,3'-Dichlorobenzidine | 1330 | 900 | 68 | 28-110 | |
| 3-Nitroaniline | 667 | 497 | 75 | 44-110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 716 | 54 | 10-110 | |
| 4-Bromophenyl phenyl ether | 667 | 571 | 86 | 39-110 | |
| 4-Chloro-3-methylphenol | 667 | 571 | 86 | 48-110 | |
| 4-Chloroaniline | 667 | 395 | 59 | 30-110 | |
| 4-Chlorophenyl phenyl ether | 667 | 549 | 82 | 40-110 | |
| 4-Nitroaniline | 667 | 588 | 88 | 48-110 | |
| 4-Nitrophenol | 1330 | 1060 | 80 | 28-110 | |
| Acenaphthene | 667 | 533 | 80 | 38-110 | |
| Acenaphthylene | 667 | 532 | 80 | 40-110 | |
| Acetophenone | 667 | 551 | 83 | 40-110 | |
| Anthracene | 667 | 545 | 82 | 48-110 | |
| Atrazine | 1330 | 1240 | 93 | 54-120 | |
| Benzaldehyde | 1330 | 1150 | 86 | 32-110 | |
| Benzo[a]anthracene | 667 | 553 | 83 | 50-110 | |
| Benzo[a]pyrene | 667 | 592 | 89 | 44-110 | |
| Benzo[b]fluoranthene | 667 | 592 | 89 | 43-110 | |
| Benzo[g,h,i]perylene | 667 | 573 | 86 | 51-110 | |
| Benzo[k]fluoranthene | 667 | 571 | 86 | 38-105 | |
| Bis(2-chloroethoxy)methane | 667 | 537 | 80 | 32-110 | |
| Bis(2-chloroethyl)ether | 667 | 533 | 80 | 34-110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 662 | 99 | 50-110 | |
| Butyl benzyl phthalate | 667 | 604 | 91 | 51-110 | |
| Caprolactam | 1330 | 1110 | 83 | 44-114 | |
| Carbazole | 667 | 564 | 85 | 50-110 | |
| Chrysene | 667 | 550 | 83 | 50-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111005.D

Lab ID: LCS 240-205786/24-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 667 | 620 | 93 | 51-110 | |
| Dibenzofuran | 667 | 548 | 82 | 43-110 | |
| Diethyl phthalate | 667 | 618 | 93 | 52-110 | |
| Dimethyl phthalate | 667 | 572 | 86 | 50-110 | |
| Di-n-butyl phthalate | 667 | 583 | 88 | 51-110 | |
| Di-n-octyl phthalate | 667 | 628 | 94 | 48-110 | |
| Fluoranthene | 667 | 539 | 81 | 51-110 | |
| Fluorene | 667 | 550 | 83 | 46-110 | |
| Hexachlorobenzene | 667 | 553 | 83 | 43-110 | |
| Hexachlorobutadiene | 667 | 559 | 84 | 29-110 | |
| Hexachlorocyclopentadiene | 667 | 240 J | 36 | 12-110 | |
| Hexachloroethane | 667 | 539 | 81 | 30-110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 606 | 91 | 50-110 | |
| Isophorone | 667 | 547 | 82 | 36-110 | |
| N-Nitrosodi-n-propylamine | 667 | 563 | 84 | 38-110 | |
| N-Nitrosodiphenylamine | 1330 | 1180 | 89 | 46-110 | |
| Naphthalene | 667 | 537 | 81 | 36-110 | |
| Nitrobenzene | 667 | 550 | 83 | 32-110 | |
| Pentachlorophenol | 1330 | 584 | 44 | 10-110 | |
| Phenanthrene | 667 | 547 | 82 | 49-110 | |
| Phenol | 667 | 542 | 81 | 38-110 | |
| Pyrene | 667 | 574 | 86 | 49-110 | |
| 3 & 4 Methylphenol | 667 | 526 | 79 | 40-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM IV
GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab File ID: 51111006.D  Lab Sample ID: MB 240-205866/8-A

Matrix: Solid  Date Extracted: 11/09/2015 12:23

Instrument ID: A4HP9  Date Analyzed: 11/11/2015 09:54

Level:(Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|---|---|---|---|
| | LCS 240-205866/9-A | 51111007.D | 11/11/2015 10:19 |
| F-VB-UST-11-14 | 240-57597-11 | 51111019.D | 11/11/2015 14:50 |
| F-VB-UST-12-14 | 240-57597-12 | 51111020.D | 11/11/2015 15:14 |
| F-VB-UST-13-14 | 240-57597-13 | 51111021.D | 11/11/2015 15:39 |
| F-VB-UST-14-14 | 240-57597-14 | 51111022.D | 11/11/2015 16:03 |
| F-VB-UST-DUP | 240-57597-15 | 51111023.D | 11/11/2015 16:28 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 51111024.D | 11/11/2015 16:52 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 51111025.D | 11/11/2015 17:17 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205866/8-A

Matrix: Solid | Lab File ID: 51111006.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 09:54

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 92-52-4 | 1,1'-Biphenyl | 3.5 | U | 50 | 3.5 |
| 108-60-1 | bis (2-chloroisopropyl) ether | 9.5 | U | 100 | 9.5 |
| 95-95-4 | 2,4,5-Trichlorophenol | 25 | U | 150 | 25 |
| 88-06-2 | 2,4,6-Trichlorophenol | 8.9 | U | 150 | 8.9 |
| 120-83-2 | 2,4-Dichlorophenol | 20 | U | 150 | 20 |
| 105-67-9 | 2,4-Dimethylphenol | 20 | U | 150 | 20 |
| 51-28-5 | 2,4-Dinitrophenol | 21 | U | 330 | 21 |
| 121-14-2 | 2,4-Dinitrotoluene | 17 | U | 200 | 17 |
| 606-20-2 | 2,6-Dinitrotoluene | 21 | U | 200 | 21 |
| 91-58-7 | 2-Chloronaphthalene | 0.45 | U | 50 | 0.45 |
| 95-57-8 | 2-Chlorophenol | 8.2 | U | 50 | 8.2 |
| 91-57-6 | 2-Methylnaphthalene | 0.50 | U | 6.7 | 0.50 |
| 95-48-7 | 2-Methylphenol | 11 | U | 200 | 11 |
| 88-74-4 | 2-Nitroaniline | 9.1 | U | 200 | 9.1 |
| 88-75-5 | 2-Nitrophenol | 8.3 | U | 50 | 8.3 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 18 | U | 100 | 18 |
| 99-09-2 | 3-Nitroaniline | 16 | U | 200 | 16 |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 9.2 | U | 150 | 9.2 |
| 101-55-3 | 4-Bromophenyl phenyl ether | 13 | U | 50 | 13 |
| 59-50-7 | 4-Chloro-3-methylphenol | 21 | U | 150 | 21 |
| 106-47-8 | 4-Chloroaniline | 17 | U | 150 | 17 |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | 13 | U | 50 | 13 |
| 100-01-6 | 4-Nitroaniline | 26 | U | 200 | 26 |
| 100-02-7 | 4-Nitrophenol | 17 | U | 330 | 17 |
| 83-32-9 | Acenaphthene | 0.76 | U | 6.7 | 0.76 |
| 208-96-8 | Acenaphthylene | 0.35 | U | 6.7 | 0.35 |
| 98-86-2 | Acetophenone | 9.2 | U | 100 | 9.2 |
| 120-12-7 | Anthracene | 0.78 | U | 6.7 | 0.78 |
| 1912-24-9 | Atrazine | 9.1 | U | 200 | 9.1 |
| 100-52-7 | Benzaldehyde | 12 | U | 100 | 12 |
| 56-55-3 | Benzo[a]anthracene | 0.63 | U | 6.7 | 0.63 |
| 50-32-8 | Benzo[a]pyrene | 0.64 | U | 6.7 | 0.64 |
| 205-99-2 | Benzo[b]fluoranthene | 0.59 | U | 6.7 | 0.59 |
| 191-24-2 | Benzo[g,h,i]perylene | 0.35 | U | 6.7 | 0.35 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205866/8-A

Matrix: Solid | Lab File ID: 51111006.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 09:54

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 207-08-9 | Benzo[k]fluoranthene | 0.68 | U | 6.7 | 0.68 |
| 111-91-1 | Bis(2-chloroethoxy)methane | 22 | U | 100 | 22 |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.0 | U | 100 | 2.0 |
| 117-81-7 | Bis(2-ethylhexyl) phthalate | 19 | U | 70 | 19 |
| 85-68-7 | Butyl benzyl phthalate | 10 | U | 70 | 10 |
| 105-60-2 | Caprolactam | 37 | U | 330 | 37 |
| 86-74-8 | Carbazole | 27 | U | 50 | 27 |
| 218-01-9 | Chrysene | 1.1 | U | 6.7 | 1.1 |
| 53-70-3 | Dibenz(a,h)anthracene | 0.66 | U | 6.7 | 0.66 |
| 132-64-9 | Dibenzofuran | 0.66 | U | 50 | 0.66 |
| 84-66-2 | Diethyl phthalate | 16 | U | 70 | 16 |
| 131-11-3 | Dimethyl phthalate | 17 | U | 70 | 17 |
| 84-74-2 | Di-n-butyl phthalate | 15 | U | 70 | 15 |
| 117-84-0 | Di-n-octyl phthalate | 7.9 | U | 70 | 7.9 |
| 206-44-0 | Fluoranthene | 0.55 | U | 6.7 | 0.55 |
| 86-73-7 | Fluorene | 0.53 | U | 6.7 | 0.53 |
| 118-74-1 | Hexachlorobenzene | 2.1 | U | 6.7 | 2.1 |
| 87-68-3 | Hexachlorobutadiene | 5.6 | U | 50 | 5.6 |
| 77-47-4 | Hexachlorocyclopentadiene | 8.1 | U | 330 | 8.1 |
| 67-72-1 | Hexachloroethane | 9.0 | U | 50 | 9.0 |
| 193-39-5 | Indeno[1,2,3-cd]pyrene | 0.35 | U | 6.7 | 0.35 |
| 78-59-1 | Isophorone | 13 | U | 50 | 13 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 6.3 | U | 50 | 6.3 |
| 86-30-6 | N-Nitrosodiphenylamine | 21 | U | 50 | 21 |
| 91-20-3 | Naphthalene | 0.82 | U | 6.7 | 0.82 |
| 98-95-3 | Nitrobenzene | 2.2 | U | 100 | 2.2 |
| 87-86-5 | Pentachlorophenol | 9.1 | U | 150 | 9.1 |
| 85-01-8 | Phenanthrene | 0.73 | U | 6.7 | 0.73 |
| 108-95-2 | Phenol | 7.3 | U | 50 | 7.3 |
| 129-00-0 | Pyrene | 0.44 | U | 6.7 | 0.44 |
| 15831-10-4 | 3 & 4 Methylphenol | 20 | U | 400 | 20 |

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205866/8-A

Matrix: Solid | Lab File ID: 51111006.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3540C | Date Extracted: 11/09/2015 12:23

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 09:54

Con. Extract Vol.: 2(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206233 | Units: ug/Kg

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 1718-51-0 | Terphenyl-d14 (Surr) | 87 | | 36-110 |
| 4165-62-2 | Phenol-d5 (Surr) | 88 | | 26-110 |
| 4165-60-0 | Nitrobenzene-d5 (Surr) | 78 | | 20-110 |
| 367-12-4 | 2-Fluorophenol (Surr) | 81 | | 24-110 |
| 321-60-8 | 2-Fluorobiphenyl (Surr) | 76 | | 24-110 |
| 118-79-6 | 2,4,6-Tribromophenol (Surr) | 49 | | 10-110 |

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111007.D

Lab ID: LCS 240-205866/9-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| 1,1'-Biphenyl | 667 | 396 | 59 | 35-110 | |
| bis (2-chloroisopropyl) ether | 667 | 514 | 77 | 29-110 | |
| 2,4,5-Trichlorophenol | 667 | 325 | 49 | 25-110 | |
| 2,4,6-Trichlorophenol | 667 | 243 | 37 | 12-110 | |
| 2,4-Dichlorophenol | 667 | 446 | 67 | 39-110 | |
| 2,4-Dimethylphenol | 667 | 444 | 67 | 29-110 | |
| 2,4-Dinitrophenol | 1330 | 464 | 35 | 10-110 | |
| 2,4-Dinitrotoluene | 667 | 473 | 71 | 48-110 | |
| 2,6-Dinitrotoluene | 667 | 462 | 69 | 45-110 | |
| 2-Chloronaphthalene | 667 | 425 | 64 | 32-110 | |
| 2-Chlorophenol | 667 | 439 | 66 | 37-110 | |
| 2-Methylnaphthalene | 667 | 428 | 64 | 36-110 | |
| 2-Methylphenol | 667 | 438 | 66 | 41-110 | |
| 2-Nitroaniline | 667 | 548 | 82 | 45-110 | |
| 2-Nitrophenol | 667 | 459 | 69 | 34-110 | |
| 3,3'-Dichlorobenzidine | 1330 | 946 | 71 | 28-110 | |
| 3-Nitroaniline | 667 | 446 | 67 | 44-110 | |
| 4,6-Dinitro-2-methylphenol | 1330 | 724 | 54 | 10-110 | |
| 4-Bromophenyl phenyl ether | 667 | 468 | 70 | 39-110 | |
| 4-Chloro-3-methylphenol | 667 | 444 | 67 | 48-110 | |
| 4-Chloroaniline | 667 | 366 | 55 | 30-110 | |
| 4-Chlorophenyl phenyl ether | 667 | 412 | 62 | 40-110 | |
| 4-Nitroaniline | 667 | 452 | 68 | 48-110 | |
| 4-Nitrophenol | 1330 | 804 | 60 | 28-110 | |
| Acenaphthene | 667 | 386 | 58 | 38-110 | |
| Acenaphthylene | 667 | 400 | 60 | 40-110 | |
| Acetophenone | 667 | 424 | 64 | 40-110 | |
| Anthracene | 667 | 462 | 69 | 48-110 | |
| Atrazine | 1330 | 1140 | 86 | 54-120 | |
| Benzaldehyde | 1330 | 1090 | 82 | 32-110 | |
| Benzo[a]anthracene | 667 | 451 | 68 | 50-110 | |
| Benzo[a]pyrene | 667 | 491 | 74 | 44-110 | |
| Benzo[b]fluoranthene | 667 | 559 | 84 | 43-110 | |
| Benzo[g,h,i]perylene | 667 | 467 | 70 | 51-110 | |
| Benzo[k]fluoranthene | 667 | 429 | 64 | 38-105 | |
| Bis(2-chloroethoxy)methane | 667 | 489 | 73 | 32-110 | |
| Bis(2-chloroethyl)ether | 667 | 503 | 75 | 34-110 | |
| Bis(2-ethylhexyl) phthalate | 667 | 544 | 82 | 50-110 | |
| Butyl benzyl phthalate | 667 | 512 | 77 | 51-110 | |
| Caprolactam | 1330 | 1000 | 75 | 44-114 | |
| Carbazole | 667 | 536 | 80 | 50-110 | |
| Chrysene | 667 | 442 | 66 | 50-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111007.D

Lab ID: LCS 240-205866/9-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 667 | 513 | 77 | 51-110 | |
| Dibenzofuran | 667 | 406 | 61 | 43-110 | |
| Diethyl phthalate | 667 | 418 | 63 | 52-110 | |
| Dimethyl phthalate | 667 | 455 | 68 | 50-110 | |
| Di-n-butyl phthalate | 667 | 538 | 81 | 51-110 | |
| Di-n-octyl phthalate | 667 | 448 | 67 | 48-110 | |
| Fluoranthene | 667 | 500 | 75 | 51-110 | |
| Fluorene | 667 | 426 | 64 | 46-110 | |
| Hexachlorobenzene | 667 | 496 | 74 | 43-110 | |
| Hexachlorobutadiene | 667 | 392 | 59 | 29-110 | |
| Hexachlorocyclopentadiene | 667 | 244 J | 37 | 12-110 | |
| Hexachloroethane | 667 | 423 | 63 | 30-110 | |
| Indeno[1,2,3-cd]pyrene | 667 | 505 | 76 | 50-110 | |
| Isophorone | 667 | 513 | 77 | 36-110 | |
| N-Nitrosodi-n-propylamine | 667 | 592 | 89 | 38-110 | |
| N-Nitrosodiphenylamine | 1330 | 987 | 74 | 46-110 | |
| Naphthalene | 667 | 387 | 58 | 36-110 | |
| Nitrobenzene | 667 | 506 | 76 | 32-110 | |
| Pentachlorophenol | 1330 | 573 | 43 | 10-110 | |
| Phenanthrene | 667 | 439 | 66 | 49-110 | |
| Phenol | 667 | 537 | 81 | 38-110 | |
| Pyrene | 667 | 481 | 72 | 49-110 | |
| 3 & 4 Methylphenol | 667 | 432 | 65 | 40-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111024.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 778 | 4.1 U | 511 | 66 | 32-110 | |
| bis (2-chloroisopropyl) ether | 778 | 11 U | 646 | 83 | 11-110 | |
| 2,4,5-Trichlorophenol | 778 | 29 U | 290 | 37 | 10-117 | |
| 2,4,6-Trichlorophenol | 778 | 10 U | 318 | 41 | 10-110 | |
| 2,4-Dichlorophenol | 778 | 23 U | 444 | 57 | 10-110 | |
| 2,4-Dimethylphenol | 778 | 23 U | 275 | 35 | 10-110 | |
| 2,4-Dinitrophenol | 1560 | 24 U | 940 | 60 | 10-110 | |
| 2,4-Dinitrotoluene | 778 | 20 U | 616 | 79 | 32-110 | |
| 2,6-Dinitrotoluene | 778 | 24 U | 597 | 77 | 35-110 | |
| 2-Chloronaphthalene | 778 | 0.52 U | 546 | 70 | 28-110 | |
| 2-Chlorophenol | 778 | 9.6 U | 504 | 65 | 10-110 | |
| 2-Methylnaphthalene | 778 | 0.58 U | 528 | 68 | 10-133 | |
| 2-Methylphenol | 778 | 13 U | 391 | 50 | 24-110 | |
| 2-Nitroaniline | 778 | 11 U | 709 | 91 | 39-110 | |
| 2-Nitrophenol | 778 | 9.7 U | 593 | 76 | 10-110 | |
| 3,3'-Dichlorobenzidine | 1560 | 21 U | 1180 | 76 | 10-110 | |
| 3-Nitroaniline | 778 | 19 U | 533 | 68 | 10-110 | |
| 4,6-Dinitro-2-methylphenol | 1560 | 11 U | 1080 | 69 | 10-110 | |
| 4-Bromophenyl phenyl ether | 778 | 15 U | 592 | 76 | 33-110 | |
| 4-Chloro-3-methylphenol | 778 | 24 U | 231 | 30 | 25-110 | |
| 4-Chloroaniline | 778 | 20 U | 391 | 50 | 10-110 | |
| 4-Chlorophenyl phenyl ether | 778 | 15 U | 537 | 69 | 32-110 | |
| 4-Nitroaniline | 778 | 30 U | 570 | 73 | 10-110 | |
| 4-Nitrophenol | 1560 | 20 U | 383 J | 25 | 10-113 | |
| Acenaphthene | 778 | 0.89 U | 501 | 64 | 22-110 | |
| Acenaphthylene | 778 | 0.41 U | 518 | 67 | 24-110 | |
| Acetophenone | 778 | 11 U | 542 | 70 | 31-110 | |
| Anthracene | 778 | 0.91 U | 585 | 75 | 20-110 | |
| Atrazine | 1560 | 11 U | 1450 | 93 | 45-118 | |
| Benzaldehyde | 1560 | 14 U | 1350 | 87 | 23-110 | |
| Benzo[a]anthracene | 778 | 0.73 U | 556 | 71 | 10-122 | |
| Benzo[a]pyrene | 778 | 0.75 U | 623 | 80 | 10-110 | |
| Benzo[b]fluoranthene | 778 | 0.69 U | 673 | 87 | 12-118 | |
| Benzo[g,h,i]perylene | 778 | 0.41 U | 598 | 77 | 10-117 | |
| Benzo[k]fluoranthene | 778 | 0.79 U | 562 | 72 | 10-121 | |
| Bis(2-chloroethoxy)methane | 778 | 26 U | 611 | 78 | 26-110 | |
| Bis(2-chloroethyl)ether | 778 | 2.3 U | 607 | 78 | 21-110 | |
| Bis(2-ethylhexyl) phthalate | 778 | 22 J | 679 | 87 | 40-110 | |
| Butyl benzyl phthalate | 778 | 12 U | 649 | 83 | 44-110 | |
| Caprolactam | 1560 | 43 U | 1260 | 81 | 10-134 | |
| Carbazole | 778 | 31 U | 666 | 86 | 34-110 | |
| Chrysene | 778 | 1.3 U | 544 | 70 | 10-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111024.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 778 | 0.77 U | 637 | 82 | 14-113 | |
| Dibenzofuran | 778 | 0.77 U | 536 | 69 | 29-110 | |
| Diethyl phthalate | 778 | 19 U | 546 | 70 | 42-110 | |
| Dimethyl phthalate | 778 | 20 U | 595 | 76 | 41-110 | |
| Di-n-butyl phthalate | 778 | 17 U | 681 | 88 | 43-110 | |
| Di-n-octyl phthalate | 778 | 9.2 U | 559 | 72 | 24-119 | |
| Fluoranthene | 778 | 0.64 U | 621 | 80 | 10-110 | |
| Fluorene | 778 | 0.62 U | 548 | 70 | 23-110 | |
| Hexachlorobenzene | 778 | 2.4 U | 623 | 80 | 34-110 | |
| Hexachlorobutadiene | 778 | 6.5 U | 472 | 61 | 25-110 | |
| Hexachlorocyclopentadiene | 778 | 9.4 U | 219 J | 28 | 10-110 | |
| Hexachloroethane | 778 | 10 U | 526 | 68 | 12-110 | |
| Indeno[1,2,3-cd]pyrene | 778 | 0.41 U | 622 | 80 | 10-114 | |
| Isophorone | 778 | 15 U | 651 | 84 | 29-110 | |
| N-Nitrosodi-n-propylamine | 778 | 7.3 U | 757 | 97 | 26-110 | |
| N-Nitrosodiphenylamine | 1560 | 24 U | 1270 | 81 | 22-110 | |
| Naphthalene | 778 | 0.96 U | 473 | 61 | 10-111 | |
| Nitrobenzene | 778 | 2.6 U | 633 | 81 | 23-110 | |
| Pentachlorophenol | 1560 | 11 U | 423 | 27 | 10-110 | |
| Phenanthrene | 778 | 0.85 U | 557 | 72 | 10-166 | |
| Phenol | 778 | 8.5 U | 395 | 51 | 17-110 | |
| Pyrene | 778 | 0.51 U | 603 | 77 | 10-147 | |
| 3 & 4 Methylphenol | 778 | 23 U | 368 J | 47 | 25-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: 51111025.D

Lab ID: 240-57597-15 MSD Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | 781 | 419 | 54 | 20 | 32 | 32-110 | |
| bis (2-chloroisopropyl) ether | 781 | 536 | 69 | 19 | 42 | 11-110 | |
| 2,4,5-Trichlorophenol | 781 | 208 | 27 | 33 | 99 | 10-117 | |
| 2,4,6-Trichlorophenol | 781 | 307 | 39 | 4 | 38 | 10-110 | |
| 2,4-Dichlorophenol | 781 | 368 | 47 | 19 | 34 | 10-110 | |
| 2,4-Dimethylphenol | 781 | 286 | 37 | 4 | 31 | 10-110 | |
| 2,4-Dinitrophenol | 1560 | 819 | 52 | 14 | 99 | 10-110 | |
| 2,4-Dinitrotoluene | 781 | 589 | 75 | 4 | 30 | 32-110 | |
| 2,6-Dinitrotoluene | 781 | 546 | 70 | 9 | 30 | 35-110 | |
| 2-Chloronaphthalene | 781 | 455 | 58 | 18 | 30 | 28-110 | |
| 2-Chlorophenol | 781 | 422 | 54 | 18 | 47 | 10-110 | |
| 2-Methylnaphthalene | 781 | 437 | 56 | 19 | 42 | 10-133 | |
| 2-Methylphenol | 781 | 416 | 53 | 6 | 51 | 24-110 | |
| 2-Nitroaniline | 781 | 646 | 83 | 9 | 31 | 39-110 | |
| 2-Nitrophenol | 781 | 492 | 63 | 19 | 49 | 10-110 | |
| 3,3'-Dichlorobenzidine | 1560 | 1210 | 77 | 2 | 56 | 10-110 | |
| 3-Nitroaniline | 781 | 551 | 71 | 3 | 30 | 10-110 | |
| 4,6-Dinitro-2-methylphenol | 1560 | 996 | 64 | 8 | 55 | 10-110 | |
| 4-Bromophenyl phenyl ether | 781 | 532 | 68 | 11 | 30 | 33-110 | |
| 4-Chloro-3-methylphenol | 781 | 272 | 35 | 16 | 54 | 25-110 | |
| 4-Chloroaniline | 781 | 392 | 50 | 0 | 36 | 10-110 | |
| 4-Chlorophenyl phenyl ether | 781 | 476 | 61 | 12 | 30 | 32-110 | |
| 4-Nitroaniline | 781 | 541 | 69 | 5 | 48 | 10-110 | |
| 4-Nitrophenol | 1560 | 446 | 29 | 15 | 49 | 10-113 | |
| Acenaphthene | 781 | 432 | 55 | 15 | 99 | 22-110 | |
| Acenaphthylene | 781 | 443 | 57 | 16 | 99 | 24-110 | |
| Acetophenone | 781 | 442 | 57 | 20 | 43 | 31-110 | |
| Anthracene | 781 | 523 | 67 | 11 | 99 | 20-110 | |
| Atrazine | 1560 | 1300 | 83 | 11 | 30 | 45-118 | |
| Benzaldehyde | 1560 | 1100 | 71 | 20 | 42 | 23-110 | |
| Benzo[a]anthracene | 781 | 536 | 69 | 4 | 99 | 10-122 | |
| Benzo[a]pyrene | 781 | 563 | 72 | 10 | 99 | 10-110 | |
| Benzo[b]fluoranthene | 781 | 638 | 82 | 5 | 99 | 12-118 | |
| Benzo[g,h,i]perylene | 781 | 545 | 70 | 9 | 99 | 10-117 | |
| Benzo[k]fluoranthene | 781 | 486 | 62 | 14 | 99 | 10-121 | |
| Bis(2-chloroethoxy)methane | 781 | 497 | 64 | 21 | 37 | 26-110 | |
| Bis(2-chloroethyl)ether | 781 | 528 | 68 | 14 | 55 | 21-110 | |
| Bis(2-ethylhexyl) phthalate | 781 | 661 | 85 | 3 | 30 | 40-110 | |
| Butyl benzyl phthalate | 781 | 614 | 79 | 5 | 30 | 44-110 | |
| Caprolactam | 1560 | 1170 | 75 | 8 | 32 | 10-134 | |
| Carbazole | 781 | 604 | 77 | 10 | 30 | 34-110 | |
| Chrysene | 781 | 521 | 67 | 4 | 99 | 10-125 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: 51111025.D

Lab ID: 240-57597-15 MSD  Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Dibenz(a,h)anthracene | 781 | 596 | 76 | 7 | 99 | 14-113 | |
| Dibenzofuran | 781 | 463 | 59 | 15 | 30 | 29-110 | |
| Diethyl phthalate | 781 | 507 | 65 | 7 | 30 | 42-110 | |
| Dimethyl phthalate | 781 | 543 | 69 | 9 | 30 | 41-110 | |
| Di-n-butyl phthalate | 781 | 631 | 81 | 8 | 30 | 43-110 | |
| Di-n-octyl phthalate | 781 | 517 | 66 | 8 | 30 | 24-119 | |
| Fluoranthene | 781 | 577 | 74 | 7 | 99 | 10-110 | |
| Fluorene | 781 | 486 | 62 | 12 | 99 | 23-110 | |
| Hexachlorobenzene | 781 | 557 | 71 | 11 | 30 | 34-110 | |
| Hexachlorobutadiene | 781 | 395 | 51 | 18 | 34 | 25-110 | |
| Hexachlorocyclopentadiene | 781 | 186 J | 24 | 16 | 79 | 10-110 | |
| Hexachloroethane | 781 | 411 | 53 | 25 | 50 | 12-110 | |
| Indeno[1,2,3-cd]pyrene | 781 | 578 | 74 | 7 | 99 | 10-114 | |
| Isophorone | 781 | 540 | 69 | 19 | 38 | 29-110 | |
| N-Nitrosodi-n-propylamine | 781 | 615 | 79 | 21 | 42 | 26-110 | |
| N-Nitrosodiphenylamine | 1560 | 1130 | 72 | 11 | 30 | 22-110 | |
| Naphthalene | 781 | 397 | 51 | 17 | 99 | 10-111 | |
| Nitrobenzene | 781 | 515 | 66 | 21 | 41 | 23-110 | |
| Pentachlorophenol | 1560 | 617 | 39 | 37 | 50 | 10-110 | |
| Phenanthrene | 781 | 510 | 65 | 9 | 99 | 10-166 | |
| Phenol | 781 | 418 | 53 | 6 | 53 | 17-110 | |
| Pyrene | 781 | 575 | 74 | 5 | 99 | 10-147 | |
| 3 & 4 Methylphenol | 781 | 392 J | 50 | 6 | 50 | 25-110 | |

# Column to be used to flag recovery and RPD values

FORM III 8270D

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 51002001.D DFTPP Injection Date: 10/02/2015

Instrument ID: A4HP10 DFTPP Injection Time: 10:22

Analysis Batch No.: 200091

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 41.5 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0)1 |
| 69 | Mass 69 relative abundance | 35.2 | |
| 70 | Less than 2.0% of mass 69 | 0.2 | (0.7)1 |
| 127 | 25.0 - 75.0% of mass 198 | 49.0 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.8 | |
| 275 | 10.0- 30.0% of mass 198 | 25.4 | |
| 365 | Greater than 0.75% of mass 198 | 3.2 | |
| 441 | Present, but less than mass 443 | 16.6 | |
| 442 | 40.0 - 110.0% of mass 198 | 102.9 | |
| 443 | 15.0 - 24.0% of mass 442 | 20.5 | (20.0)2 |

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD5 240-200091/2 | 51002002.D | 10/02/2015 | 10:48 |
| | STD4 240-200091/3 | 51002003.D | 10/02/2015 | 11:12 |
| | STD3 240-200091/4 | 51002004.D | 10/02/2015 | 11:37 |
| | STD2 240-200091/5 | 51002005.D | 10/02/2015 | 12:01 |
| | STD1 240-200091/6 | 51002006.D | 10/02/2015 | 12:25 |
| | STD9 240-200091/7 | 51002007.D | 10/02/2015 | 12:50 |
| | STD8 240-200091/8 | 51002008.D | 10/02/2015 | 13:15 |
| | STD7 240-200091/9 | 51002009.D | 10/02/2015 | 13:39 |
| | STD6 240-200091/10 | 51002010.D | 10/02/2015 | 14:04 |
| | ICV 240-200091/11 | 51002011.D | 10/02/2015 | 14:28 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48  Calibration End Date: 10/02/2015 14:04  Calibration ID: 30980

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-200091/6 | 51002006.D |
| Level 2 | STD2 240-200091/5 | 51002005.D |
| Level 3 | STD3 240-200091/4 | 51002004.D |
| Level 4 | STD4 240-200091/3 | 51002003.D |
| Level 5 | STD5 240-200091/2 | 51002002.D |
| Level 6 | STD6 240-200091/10 | 51002010.D |
| Level 7 | STD7 240-200091/9 | 51002009.D |
| Level 8 | STD8 240-200091/8 | 51002008.D |
| Level 9 | STD9 240-200091/7 | 51002007.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,4-Dioxane | 0.6695 | 0.7664<br>0.6450 | 0.7112<br>0.6601 | 0.6791<br>0.6797 | 0.7113 | Ave | | 0.6903 | | | | 5.6 | | 20.0 | | | |
| N-Nitrosodimethylamine | 0.9213 | 0.9861<br>0.9036 | 0.9021<br>0.9073 | 0.9106<br>0.9353 | 0.9364 | Ave | | 0.9253 | | | | 3.0 | | 20.0 | | | |
| Pyridine | 1.6752 | 1.7590<br>1.6375 | 1.6236<br>1.6765 | 1.6397<br>1.7157 | 1.7460 | Ave | | 1.6842 | | | | 3.0 | | 20.0 | | | |
| Benzaldehyde | 1.1161 | 1.2599<br>0.9941 | 1.2278<br>0.9643 | 1.2174<br>0.9002 | 1.2082 | Ave | | 1.1110 | | | 0.0100 | 12.6 | | 20.0 | | | |
| Phenol | 1.8772 | 2.0486<br>1.7989 | 1.9063<br>1.8076 | 1.9409<br>1.8084 | 1.9432 | Ave | | 1.8914 | | | 0.8000 | 4.6 | | 20.0 | | | |
| Aniline | 2.4273 | 2.4891<br>2.3113 | 2.4030<br>2.3297 | 2.4066<br>2.3780 | 2.4676 | Ave | | 2.4016 | | | | 2.6 | | 20.0 | | | |
| Bis(2-chloroethyl)ether | 1.5242 | 1.5993<br>1.4438 | 1.5227<br>1.4462 | 1.5177<br>1.4436 | 1.5534 | Ave | | 1.5064 | | | 0.7000 | 3.8 | | 20.0 | | | |
| 2-Chlorophenol | 1.5436 | 1.5843<br>1.4930 | 1.5903<br>1.5124 | 1.5578<br>1.5091 | 1.5573 | Ave | | 1.5435 | | | 0.8000 | 2.3 | | 20.0 | | | |
| n-Decane | 1.8806 | 2.1172<br>1.7823 | 2.0495<br>1.7515 | 1.9665<br>1.7279 | 2.0152 | Ave | | 1.9113 | | | | 7.7 | | 20.0 | | | |
| 1,3-Dichlorobenzene | 1.5929 | 1.6788<br>1.5009 | 1.5686<br>1.5104 | 1.6231<br>1.5138 | 1.6452 | Ave | | 1.5792 | | | | 4.3 | | 20.0 | | | |
| 1,4-Dichlorobenzene | 1.6286 | 1.7289<br>1.5274 | 1.6483<br>1.5572 | 1.6056<br>1.5444 | 1.6767 | Ave | | 1.6146 | | | | 4.3 | | 20.0 | | | |
| Benzyl alcohol | 1.0249 | 1.1085<br>0.9791 | 1.0175<br>0.9932 | 1.0225<br>0.9932 | 1.0587 | Ave | | 1.0247 | | | | 4.1 | | 20.0 | | | |
| 1,2-Dichlorobenzene | 1.5382 | 1.6271<br>1.4641 | 1.5700<br>1.4793 | 1.5350<br>1.4835 | 1.5516 | Ave | | 1.5311 | | | | 3.5 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48  Calibration End Date: 10/02/2015 14:04  Calibration ID: 30980

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2-Methylphenol | 1.4114 | 1.4902<br>1.3491 | 1.4012<br>1.3603 | 1.4223<br>1.3516 | 1.5104 | Ave | | 1.4121 | | | 0.7000 | 4.3 | | 20.0 | | | |
| bis (2-chloroisopropyl) ether | 2.1032 | 2.4010<br>1.9023 | 2.3063<br>1.8708 | 2.2908<br>1.8368 | 2.2583 | Ave | | 2.1212 | | | | 10.6 | | 20.0 | | | |
| Indene | 2.5417 | 2.8532<br>2.3321 | 2.6823<br>2.2562 | 2.7190<br>2.1708 | 2.7151 | Ave | | 2.5338 | | | | 9.9 | | 20.0 | | | |
| 3 & 4 Methylphenol | 1.4717 | 1.5716<br>1.3921 | 1.4945<br>1.3910 | 1.4916<br>1.3913 | 1.5495 | Ave | | 1.4692 | | | | 4.9 | | 20.0 | | | |
| N-Nitrosodi-n-propylamine | 0.9027 | 0.9251<br>0.8420 | 0.9104<br>0.8191 | 0.8990<br>0.7921 | 0.9376 | Ave | | 0.8785 | | | 0.5000 | 6.1 | | 20.0 | | | |
| Acetophenone | 2.0138 | 2.0766<br>1.8967 | 2.0491<br>1.8871 | 2.0752<br>1.8264 | 2.0787 | Ave | | 1.9880 | | | 0.0100 | 5.1 | | 20.0 | | | |
| Hexachloroethane | 0.6506 | 0.6511<br>0.6153 | 0.6147<br>0.6272 | 0.6386<br>0.6127 | 0.6770 | Ave | | 0.6359 | | | 0.3000 | 3.6 | | 20.0 | | | |
| Nitrobenzene | 0.3700 | 0.3890<br>0.3608 | 0.3813<br>0.3557 | 0.3686<br>0.3512 | 0.3786 | Ave | | 0.3694 | | | 0.2000 | 3.6 | | 20.0 | | | |
| Isophorone | 0.6883 | 0.7238<br>0.6711 | 0.6929<br>0.6633 | 0.6847<br>0.6524 | 0.7054 | Ave | | 0.6852 | | | 0.4000 | 3.4 | | 20.0 | | | |
| 2-Nitrophenol | 0.1876 | 0.1380<br>0.1801 | 0.1435<br>0.1805 | 0.1564<br>0.1754 | 0.1739 | Ave | | 0.1669 | | | 0.1000 | 11.1 | | 20.0 | | | |
| 2,4-Dimethylphenol | 0.3616 | 0.3894<br>0.3430 | 0.3809<br>0.3374 | 0.3682<br>0.3237 | 0.3597 | Ave | | 0.3580 | | | 0.2000 | 6.2 | | 20.0 | | | |
| Benzoic acid | 0.2291 | +++++<br>0.2141 | 0.0903<br>0.2178 | 0.1351<br>0.2228 | 0.1906 | Lin1 | -0.305 | 0.2283 | | | | | | | 0.9990 | | 0.9900 |
| Bis(2-chloroethoxy)methane | 0.4177 | 0.4647<br>0.3986 | 0.4508<br>0.3937 | 0.4375<br>0.3846 | 0.4412 | Ave | | 0.4236 | | | 0.3000 | 6.9 | | 20.0 | | | |
| 2,4-Dichlorophenol | 0.3004 | 0.3102<br>0.2890 | 0.2958<br>0.2884 | 0.2907<br>0.2810 | 0.2969 | Ave | | 0.2940 | | | 0.2000 | 3.0 | | 20.0 | | | |
| 1,2,4-Trichlorobenzene | 0.3054 | 0.3141<br>0.2927 | 0.3209<br>0.2858 | 0.3095<br>0.2778 | 0.3117 | Ave | | 0.3022 | | | | 5.0 | | 20.0 | | | |
| Naphthalene | 1.3004<br>1.0690 | 1.1308<br>1.0304 | 1.0735<br>1.0093 | 1.0684<br>0.9772 | 1.0969 | Ave | | 1.0840 | | | 0.7000 | 8.6 | | 20.0 | | | |
| 4-Chloroaniline | 0.4741 | 0.4930<br>0.4559 | 0.5069<br>0.4525 | 0.4937<br>0.4485 | 0.4800 | Ave | | 0.4756 | | | 0.0100 | 4.6 | | 20.0 | | | |
| 2,6-Dichlorophenol | 0.2958 | 0.3215<br>0.2856 | 0.3004<br>0.2833 | 0.2958<br>0.2781 | 0.2986 | Ave | | 0.2949 | | | | 4.5 | | 20.0 | | | |
| Hexachlorobutadiene | 0.1890 | 0.2046<br>0.1802 | 0.1919<br>0.1770 | 0.1878<br>0.1720 | 0.1923 | Ave | | 0.1868 | | | 0.0100 | 5.5 | | 20.0 | | | |
| Caprolactam | 0.1158 | 0.1186<br>0.1145 | 0.1187<br>0.1160 | 0.1123<br>0.1173 | 0.1180 | Ave | | 0.1164 | | | 0.0100 | 1.9 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48  Calibration End Date: 10/02/2015 14:04  Calibration ID: 30980

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 4-Chloro-3-methylphenol | 0.3081 | 0.3124<br>0.2963 | 0.3078<br>0.2958 | 0.3032<br>0.2895 | 0.3130 | Ave | | 0.3033 | | | 0.2000 | 2.8 | | 20.0 | | | |
| 2-Methylnaphthalene | 0.7756<br>0.6955 | 0.7490<br>0.6693 | 0.7170<br>0.6556 | 0.6957<br>0.6313 | 0.7165 | Ave | | 0.7006 | | | 0.4000 | 6.5 | | 20.0 | | | |
| 1-Methylnaphthalene | 0.7676<br>0.6107 | 0.6360<br>0.5848 | 0.6456<br>0.5768 | 0.6162<br>0.5626 | 0.6282 | Ave | | 0.6254 | | | | 9.6 | | 20.0 | | | |
| Hexachlorocyclopentadiene | 0.2948 | 0.2072<br>0.2918 | 0.2317<br>0.2918 | 0.2528<br>0.2873 | 0.2937 | Ave | | 0.2689 | | | 0.0500 | 12.7 | | 20.0 | | | |
| 1,2,4,5-Tetrachlorobenzene | 0.4885 | 0.5652<br>0.4629 | 0.5402<br>0.4463 | 0.5472<br>0.4297 | 0.5183 | Ave | | 0.4998 | | | 0.0100 | 10.1 | | 20.0 | | | |
| 2,4,6-Trichlorophenol | 0.3715 | 0.3724<br>0.3588 | 0.3515<br>0.3509 | 0.3712<br>0.3425 | 0.3759 | Ave | | 0.3619 | | | 0.2000 | 3.5 | | 20.0 | | | |
| 2,4,5-Trichlorophenol | 0.3894 | 0.3736<br>0.3765 | 0.3808<br>0.3754 | 0.3750<br>0.3713 | 0.3936 | Ave | | 0.3794 | | | 0.2000 | 2.1 | | 20.0 | | | |
| 1,1'-Biphenyl | 1.6748<br>1.3725 | 1.4785<br>1.3090 | 1.4844<br>1.2882 | 1.4329<br>1.2449 | 1.3992 | Ave | | 1.4094 | | | 0.0100 | 9.2 | | 20.0 | | | |
| 2-Chloronaphthalene | 1.4430<br>1.1165 | 1.3094<br>1.0904 | 1.1919<br>1.0799 | 1.2036<br>1.0476 | 1.1544 | Ave | | 1.1819 | | | 0.8000 | 10.7 | | 20.0 | | | |
| 2-Nitroaniline | 0.3659 | 0.3458<br>0.3517 | 0.3335<br>0.3522 | 0.3394<br>0.3426 | 0.3714 | Ave | | 0.3503 | | | 0.0100 | 3.7 | | 20.0 | | | |
| Dimethyl phthalate | 1.2918 | 1.3757<br>1.2433 | 1.3134<br>1.2373 | 1.2978<br>1.2027 | 1.3157 | Ave | | 1.2847 | | | 0.0100 | 4.3 | | 20.0 | | | |
| 1,3-Dinitrobenzene | 0.2128 | 0.1253<br>0.2137 | 0.1311<br>0.2169 | 0.1611<br>0.2086 | 0.1876 | Lin1 | -0.066 | 0.2149 | | | | | | | 0.9980 | | 0.9900 |
| 2,6-Dinitrotoluene | 0.2980 | 0.2441<br>0.2879 | 0.2649<br>0.2943 | 0.2864<br>0.2823 | 0.2962 | Ave | | 0.2818 | | | | 6.6 | | 20.0 | | | |
| Acenaphthylene | 2.0743<br>1.7077 | 1.8368<br>1.6563 | 1.7806<br>1.6705 | 1.7896<br>1.5999 | 1.7512 | Ave | | 1.7630 | | | 0.9000 | 7.8 | | 20.0 | | | |
| 3-Nitroaniline | 0.3380 | 0.2955<br>0.3324 | 0.3068<br>0.3328 | 0.3220<br>0.3280 | 0.3359 | Ave | | 0.3239 | | | 0.0100 | 4.7 | | 20.0 | | | |
| 2,4-Dinitrophenol | 0.1481 | +++++<br>0.1538 | 0.0351<br>0.1544 | 0.0510<br>0.1462 | 0.0918 | Lin1 | -0.328 | 0.1574 | | | 0.0100 | | | | 0.9920 | | 0.9900 |
| Acenaphthene | 1.3039<br>1.0225 | 1.2131<br>0.9782 | 1.1513<br>0.9499 | 1.1017<br>0.8846 | 1.0737 | Ave | | 1.0754 | | | 0.9000 | 12.4 | | 20.0 | | | |
| 4-Nitrophenol | 0.1794 | 0.1547<br>0.1768 | 0.1585<br>0.1774 | 0.1663<br>0.1698 | 0.1787 | Ave | | 0.1702 | | | | 5.6 | | 20.0 | | | |
| 2,4-Dinitrotoluene | 0.4036 | 0.2963<br>0.3894 | 0.3270<br>0.3928 | 0.3557<br>0.3847 | 0.3830 | Ave | | 0.3666 | | | 0.2000 | 10.2 | | 20.0 | | | |
| Dibenzofuran | 1.9069<br>1.5721 | 1.7236<br>1.5073 | 1.6855<br>1.5083 | 1.6327<br>1.4681 | 1.6051 | Ave | | 1.6233 | | | 0.8000 | 8.4 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48  Calibration End Date: 10/02/2015 14:04  Calibration ID: 30980

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2,3,4,6-Tetrachlorophenol | 0.3261 | 0.3032<br>0.3202 | 0.3101<br>0.3234 | 0.3209<br>0.3127 | 0.3268 | Ave | | 0.3179 | | | 0.0100 | 2.6 | | 20.0 | | | |
| Diethyl phthalate | 1.1755 | 1.3606<br>1.0860 | 1.2751<br>1.0391 | 1.2561<br>0.9672 | 1.2198 | Ave | | 1.1724 | | | 0.0100 | 11.3 | | 20.0 | | | |
| Hexadecane | 0.5384 | 0.6539<br>0.4892 | 0.6552<br>0.4399 | 0.6420<br>0.3964 | 0.6215 | Ave | | 0.5546 | | | | 18.6 | | 20.0 | | | |
| 4-Chlorophenyl phenyl ether | 0.6018 | 0.6868<br>0.5811 | 0.6476<br>0.5764 | 0.6229<br>0.5609 | 0.6109 | Ave | | 0.6110 | | | 0.4000 | 6.8 | | 20.0 | | | |
| 4-Nitroaniline | 0.3283 | 0.3168<br>0.3054 | 0.3309<br>0.2952 | 0.3259<br>0.2745 | 0.3421 | Ave | | 0.3149 | | | 0.0100 | 7.0 | | 20.0 | | | |
| Fluorene | 1.5331<br>1.3077 | 1.4393<br>1.2410 | 1.3917<br>1.1745 | 1.3621<br>1.1146 | 1.3380 | Ave | | 1.3224 | | | 0.9000 | 9.9 | | 20.0 | | | |
| 4,6-Dinitro-2-methylphenol | 0.1405 | +++++<br>0.1420 | 0.0535<br>0.1473 | 0.0736<br>0.1406 | 0.1077 | Lin1 | -0.238 | 0.1483 | | | 0.0100 | | | | 0.9970 | | 0.9900 |
| N-Nitrosodiphenylamine | 0.5721 | 0.6386<br>0.5413 | 0.6032<br>0.5217 | 0.6014<br>0.4871 | 0.6027 | Ave | | 0.5710 | | | 0.0100 | 8.9 | | 20.0 | | | |
| 1,2-Diphenylhydrazine(as Azobenzene) | 0.7979 | 0.8327<br>0.7726 | 0.8253<br>0.7603 | 0.8422<br>0.7222 | 0.8418 | Ave | | 0.7997 | | | | 5.5 | | 20.0 | | | |
| Azobenzene | 0.7979 | 0.8316<br>0.7731 | 0.8275<br>0.7602 | 0.8424<br>0.7228 | 0.8424 | Ave | | 0.7997 | | | | 5.5 | | 20.0 | | | |
| 4-Bromophenyl phenyl ether | 0.2291 | 0.2468<br>0.2220 | 0.2422<br>0.2182 | 0.2340<br>0.2104 | 0.2360 | Ave | | 0.2298 | | | 0.1000 | 5.4 | | 20.0 | | | |
| Hexachlorobenzene | 0.2897<br>0.2474 | 0.2906<br>0.2387 | 0.2639<br>0.2357 | 0.2702<br>0.2283 | 0.2592 | Ave | | 0.2582 | | | 0.1000 | 8.8 | | 20.0 | | | |
| Atrazine | 0.2085 | 0.2470<br>0.1989 | 0.2352<br>0.1914 | 0.2302<br>0.1784 | 0.2226 | Ave | | 0.2140 | | | 0.0100 | 11.1 | | 20.0 | | | |
| Pentachlorophenol | 0.1566 | 0.1203<br>0.1461 | 0.1373<br>0.1435 | 0.1462<br>0.1377 | 0.1580 | Ave | | 0.1432 | | | 0.0500 | 8.4 | | 20.0 | | | |
| n-Octadecane | 0.7811 | 1.0785<br>0.6664 | 1.0384<br>0.6036 | 1.0135<br>0.5462 | 0.9376 | Qua | 0.2920 | 0.8735 | -0.013823 | | | | | | 0.9980 | | 0.9900 |
| Phenanthrene | 1.3748<br>1.1479 | 1.2369<br>1.1131 | 1.1926<br>1.0919 | 1.1939<br>1.0530 | 1.1921 | Ave | | 1.1773 | | | 0.7000 | 8.0 | | 20.0 | | | |
| Anthracene | 1.3205<br>1.1782 | 1.2586<br>1.1423 | 1.2160<br>1.1412 | 1.1785<br>1.0647 | 1.2105 | Ave | | 1.1901 | | | 0.7000 | 6.2 | | 20.0 | | | |
| Carbazole | 1.1072 | 1.2431<br>1.0787 | 1.2009<br>1.0743 | 1.1353<br>1.0193 | 1.1600 | Ave | | 1.1273 | | | 0.0100 | 6.5 | | 20.0 | | | |
| Di-n-butyl phthalate | 1.3947 | 1.4647<br>1.3496 | 1.4241<br>1.3237 | 1.4029<br>1.2334 | 1.4453 | Ave | | 1.3798 | | | 0.0100 | 5.5 | | 20.0 | | | |
| Fluoranthene | 1.5321<br>1.2882 | 1.3939<br>1.2469 | 1.3537<br>1.2340 | 1.3252<br>1.1725 | 1.3424 | Ave | | 1.3210 | | | 0.6000 | 7.9 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 200091

SDG No.:

Instrument ID: A4HP10  GC Column: DB-624  ID: 0.18(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/02/2015 10:48  Calibration End Date: 10/02/2015 14:04  Calibration ID: 30980

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Benzidine | 0.6451 | 0.5338<br>0.6050 | 0.5417<br>0.5945 | 0.5546<br>0.5565 | 0.6428 | Ave | | 0.5843 | | | | 7.6 | | 20.0 | | | |
| Pyrene | 1.3418<br>1.1061 | 1.1896<br>1.0765 | 1.1446<br>1.0713 | 1.1384<br>1.0412 | 1.1040 | Ave | | 1.1348 | | | 0.6000 | 7.9 | | 20.0 | | | |
| Butyl benzyl phthalate | 0.5176 | 0.5409<br>0.5064 | 0.5195<br>0.5022 | 0.5221<br>0.4846 | 0.5206 | Ave | | 0.5142 | | | 0.0100 | 3.2 | | 20.0 | | | |
| 3,3'-Dichlorobenzidine | 0.3938 | 0.4187<br>0.3806 | 0.4084<br>0.3666 | 0.4148<br>0.3572 | 0.4218 | Ave | | 0.3952 | | | 0.0100 | 6.3 | | 20.0 | | | |
| Bis(2-ethylhexyl) phthalate | 0.6563 | 0.7249<br>0.6056 | 0.7320<br>0.5577 | 0.7423<br>0.5175 | 0.7272 | Ave | | 0.6579 | | | 0.0100 | 13.4 | | 20.0 | | | |
| Benzo[a]anthracene | 1.2994<br>1.1072 | 1.1154<br>1.0880 | 1.1109<br>1.1166 | 1.1030<br>1.0814 | 1.1126 | Ave | | 1.1261 | | | 0.8000 | 5.9 | | 20.0 | | | |
| Chrysene | 1.2630<br>1.0330 | 1.1264<br>1.0316 | 1.0723<br>1.0239 | 1.0588<br>1.0141 | 1.0605 | Ave | | 1.0759 | | | 0.7000 | 7.2 | | 20.0 | | | |
| Di-n-octyl phthalate | 1.4250 | 1.2463<br>1.3943 | 1.3097<br>1.4235 | 1.3592<br>1.3602 | 1.4489 | Ave | | 1.3709 | | | 0.0100 | 4.9 | | 20.0 | | | |
| Benzo[b]fluoranthene | 1.1428<br>1.1817 | 1.1483<br>1.0957 | 1.1239<br>1.1339 | 1.0694<br>1.1422 | 1.1127 | Ave | | 1.1278 | | | 0.7000 | 2.9 | | 20.0 | | | |
| Benzo[k]fluoranthene | 1.4480<br>1.2152 | 1.3389<br>1.2296 | 1.3013<br>1.2500 | 1.2492<br>1.1684 | 1.2686 | Ave | | 1.2743 | | | 0.7000 | 6.4 | | 20.0 | | | |
| Benzo[a]pyrene | 1.1242<br>1.1335 | 1.1003<br>1.1221 | 1.1316<br>1.1430 | 1.0872<br>1.1046 | 1.1374 | Ave | | 1.1204 | | | 0.7000 | 1.7 | | 20.0 | | | |
| Indeno[1,2,3-cd]pyrene | 1.0594<br>1.2888 | 1.1682<br>1.2677 | 1.2083<br>1.2995 | 1.2340<br>1.2525 | 1.3208 | Ave | | 1.2332 | | | 0.5000 | 6.5 | | 20.0 | | | |
| Dibenz(a,h)anthracene | 0.7887<br>1.0727 | 1.0310<br>1.0681 | 1.0191<br>1.0966 | 1.0137<br>1.0546 | 1.0748 | Ave | | 1.0244 | | | 0.4000 | 9.0 | | 20.0 | | | |
| Benzo[g,h,i]perylene | 0.9436<br>1.1332 | 1.0999<br>1.1069 | 1.0654<br>1.1485 | 1.0731<br>1.1229 | 1.1388 | Ave | | 1.0925 | | | 0.5000 | 5.7 | | 20.0 | | | |
| 2-Fluorophenol (Surr) | 1.4936 | 1.4868<br>1.4489 | 1.4884<br>1.4663 | 1.5162<br>1.4709 | 1.5176 | Ave | | 1.4861 | | | | 1.6 | | 20.0 | | | |
| Phenol-d5 (Surr) | 1.8356 | 1.9501<br>1.7490 | 1.8116<br>1.7382 | 1.7830<br>1.7442 | 1.8996 | Ave | | 1.8139 | | | | 4.3 | | 20.0 | | | |
| Nitrobenzene-d5 (Surr) | 0.4177<br>0.3995 | 0.4058<br>0.3874 | 0.3992<br>0.3905 | 0.3968<br>0.3818 | 0.4054 | Ave | | 0.3982 | | | | 2.7 | | 20.0 | | | |
| 2-Fluorobiphenyl (Surr) | 1.6344<br>1.2399 | 1.4109<br>1.1877 | 1.3265<br>1.1814 | 1.2886<br>1.1377 | 1.2630 | Ave | | 1.2967 | | | | 11.7 | | 20.0 | | | |
| 2,4,6-Tribromophenol (Surr) | 0.1986 | 0.1782<br>0.1920 | 0.1845<br>0.1971 | 0.1810<br>0.1944 | 0.1873 | Ave | | 0.1891 | | | | 4.0 | | 20.0 | | | |
| Terphenyl-d14 (Surr) | 0.9059<br>0.7324 | 0.8249<br>0.7190 | 0.7668<br>0.7171 | 0.7574<br>0.6951 | 0.7388 | Ave | | 0.7619 | | | | 8.6 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-200091/11 Calibration Date: 10/02/2015 14:28

Instrument ID: A4HP10 Calib Start Date: 10/02/2015 10:48

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/02/2015 14:04

Lab File ID: 51002011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.6903 | 0.6472 | | 9.38 | 10.0 | -6.2 | 30.0 |
| N-Nitrosodimethylamine | Ave | 0.9253 | 0.7776 | | 8.40 | 10.0 | -16.0 | 30.0 |
| Pyridine | Ave | 1.684 | 1.634 | | 9.70 | 10.0 | -3.0 | 30.0 |
| Benzaldehyde | Ave | 1.111 | 1.071 | 0.0100 | 19.3 | 20.0 | -3.6 | 30.0 |
| Phenol | Ave | 1.891 | 1.832 | 0.8000 | 9.69 | 10.0 | -3.1 | 30.0 |
| Aniline | Ave | 2.402 | 2.297 | | 9.57 | 10.0 | -4.3 | 30.0 |
| Bis(2-chloroethyl)ether | Ave | 1.506 | 1.415 | 0.7000 | 9.40 | 10.0 | -6.0 | 30.0 |
| 2-Chlorophenol | Ave | 1.543 | 1.512 | 0.8000 | 9.80 | 10.0 | -2.0 | 30.0 |
| n-Decane | Ave | 1.911 | 1.813 | | 9.48 | 10.0 | -5.2 | 30.0 |
| 1,3-Dichlorobenzene | Ave | 1.579 | 1.531 | | 9.69 | 10.0 | -3.1 | 30.0 |
| 1,4-Dichlorobenzene | Ave | 1.615 | 1.563 | | 9.68 | 10.0 | -3.2 | 30.0 |
| Benzyl alcohol | Ave | 1.025 | 0.9929 | | 9.69 | 10.0 | -3.1 | 30.0 |
| 1,2-Dichlorobenzene | Ave | 1.531 | 1.463 | | 9.56 | 10.0 | -4.4 | 30.0 |
| 2-Methylphenol | Ave | 1.412 | 1.403 | 0.7000 | 9.93 | 10.0 | -0.7 | 30.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.121 | 2.000 | | 9.43 | 10.0 | -5.7 | 30.0 |
| Indene | Ave | 2.534 | 2.356 | | 18.6 | 20.0 | -7.0 | 30.0 |
| 3 & 4 Methylphenol | Ave | 1.469 | 1.437 | | 9.78 | 10.0 | -2.2 | 30.0 |
| N-Nitrosodi-n-propylamine | Ave | 0.8785 | 0.8794 | 0.5000 | 10.0 | 10.0 | 0.1 | 30.0 |
| Acetophenone | Ave | 1.988 | 1.939 | 0.0100 | 9.76 | 10.0 | -2.4 | 30.0 |
| Hexachloroethane | Ave | 0.6359 | 0.6318 | 0.3000 | 9.94 | 10.0 | -0.6 | 30.0 |
| Nitrobenzene | Ave | 0.3694 | 0.3535 | 0.2000 | 9.57 | 10.0 | -4.3 | 30.0 |
| Isophorone | Ave | 0.6852 | 0.6871 | 0.4000 | 10.0 | 10.0 | 0.3 | 30.0 |
| 2-Nitrophenol | Ave | 0.1669 | 0.1805 | 0.1000 | 10.8 | 10.0 | 8.1 | 30.0 |
| 2,4-Dimethylphenol | Ave | 0.3580 | 0.3438 | 0.2000 | 9.61 | 10.0 | -3.9 | 30.0 |
| Benzoic acid | Lin1 | | 0.2265 | | 21.2 | 20.0 | 5.9 | 30.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4236 | 0.4028 | 0.3000 | 9.51 | 10.0 | -4.9 | 30.0 |
| 2,4-Dichlorophenol | Ave | 0.2940 | 0.2826 | 0.2000 | 9.61 | 10.0 | -3.9 | 30.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.3022 | 0.2949 | | 9.76 | 10.0 | -2.4 | 30.0 |
| Naphthalene | Ave | 1.084 | 1.032 | 0.7000 | 9.52 | 10.0 | -4.8 | 30.0 |
| 4-Chloroaniline | Ave | 0.4756 | 0.4464 | 0.0100 | 9.39 | 10.0 | -6.1 | 30.0 |
| 2,6-Dichlorophenol | Ave | 0.2949 | 0.2868 | | 9.73 | 10.0 | -2.7 | 30.0 |
| Hexachlorobutadiene | Ave | 0.1868 | 0.1827 | 0.0100 | 9.78 | 10.0 | -2.2 | 30.0 |
| Caprolactam | Ave | 0.1164 | 0.1183 | 0.0100 | 20.3 | 20.0 | 1.6 | 30.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3033 | 0.2997 | 0.2000 | 9.88 | 10.0 | -1.2 | 30.0 |
| 2-Methylnaphthalene | Ave | 0.7006 | 0.6481 | 0.4000 | 9.25 | 10.0 | -7.5 | 30.0 |
| 1-Methylnaphthalene | Ave | 0.6254 | 0.6114 | | 9.78 | 10.0 | -2.2 | 30.0 |
| Hexachlorocyclopentadiene | Ave | 0.2689 | 0.2968 | 0.0500 | 11.0 | 10.0 | 10.4 | 30.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4998 | 0.4717 | 0.0100 | 9.44 | 10.0 | -5.6 | 30.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3619 | 0.3701 | 0.2000 | 10.2 | 10.0 | 2.3 | 30.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3794 | 0.3824 | 0.2000 | 10.1 | 10.0 | 0.8 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-200091/11 Calibration Date: 10/02/2015 14:28

Instrument ID: A4HP10 Calib Start Date: 10/02/2015 10:48

GC Column: DB-624 ID: 0.18(mm) Calib End Date: 10/02/2015 14:04

Lab File ID: 51002011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.409 | 1.323 | 0.0100 | 9.39 | 10.0 | -6.1 | 30.0 |
| 2-Chloronaphthalene | Ave | 1.182 | 1.109 | 0.8000 | 9.38 | 10.0 | -6.2 | 30.0 |
| 2-Nitroaniline | Ave | 0.3503 | 0.3633 | 0.0100 | 10.4 | 10.0 | 3.7 | 30.0 |
| Dimethyl phthalate | Ave | 1.285 | 1.249 | 0.0100 | 9.72 | 10.0 | -2.8 | 30.0 |
| 1,3-Dinitrobenzene | Lin1 | | 0.2089 | | 10.0 | 10.0 | 0.3 | 30.0 |
| 2,6-Dinitrotoluene | Ave | 0.2818 | 0.2952 | | 10.5 | 10.0 | 4.8 | 30.0 |
| Acenaphthylene | Ave | 1.763 | 1.642 | 0.9000 | 9.31 | 10.0 | -6.9 | 30.0 |
| 3-Nitroaniline | Ave | 0.3239 | 0.3357 | 0.0100 | 10.4 | 10.0 | 3.6 | 30.0 |
| 2,4-Dinitrophenol | Lin1 | | 0.1483 | 0.0100 | 20.9 | 20.0 | 4.7 | 30.0 |
| Acenaphthene | Ave | 1.075 | 1.012 | 0.9000 | 9.41 | 10.0 | -5.9 | 30.0 |
| 4-Nitrophenol | Ave | 0.1702 | 0.1772 | | 20.8 | 20.0 | 4.1 | 30.0 |
| 2,4-Dinitrotoluene | Ave | 0.3666 | 0.3977 | 0.2000 | 10.8 | 10.0 | 8.5 | 30.0 |
| Dibenzofuran | Ave | 1.623 | 1.525 | 0.8000 | 9.39 | 10.0 | -6.1 | 30.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.3179 | 0.3218 | 0.0100 | 10.1 | 10.0 | 1.2 | 30.0 |
| Diethyl phthalate | Ave | 1.172 | 1.149 | 0.0100 | 9.80 | 10.0 | -2.0 | 30.0 |
| Hexadecane | Ave | 0.5546 | 0.5084 | | 9.17 | 10.0 | -8.3 | 30.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.6110 | 0.5884 | 0.4000 | 9.63 | 10.0 | -3.7 | 30.0 |
| 4-Nitroaniline | Ave | 0.3149 | 0.3341 | 0.0100 | 10.6 | 10.0 | 6.1 | 30.0 |
| Fluorene | Ave | 1.322 | 1.271 | 0.9000 | 9.61 | 10.0 | -3.9 | 30.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1420 | 0.0100 | 20.8 | 20.0 | 3.8 | 30.0 |
| N-Nitrosodiphenylamine | Ave | 0.5710 | 0.5568 | 0.0100 | 19.5 | 20.0 | -2.5 | 30.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 0.7994 | 0.8227 | | 10.3 | 10.0 | 2.9 | 30.0 |
| Azobenzene | Ave | 0.7997 | 0.8228 | | 10.3 | 10.0 | 2.9 | 30.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.2298 | 0.2289 | 0.1000 | 9.96 | 10.0 | -0.4 | 30.0 |
| Hexachlorobenzene | Ave | 0.2582 | 0.2429 | 0.1000 | 9.41 | 10.0 | -5.9 | 30.0 |
| Atrazine | Ave | 0.2140 | 0.2028 | 0.0100 | 19.0 | 20.0 | -5.2 | 30.0 |
| n-Octadecane | Qua | | 0.7442 | | 9.66 | 10.0 | -3.4 | 30.0 |
| Pentachlorophenol | Ave | 0.1432 | 0.1612 | 0.0500 | 22.5 | 20.0 | 12.5 | 30.0 |
| Phenanthrene | Ave | 1.177 | 1.107 | 0.7000 | 9.41 | 10.0 | -5.9 | 30.0 |
| Anthracene | Ave | 1.190 | 1.153 | 0.7000 | 9.69 | 10.0 | -3.1 | 30.0 |
| Carbazole | Ave | 1.127 | 1.062 | 0.0100 | 9.42 | 10.0 | -5.8 | 30.0 |
| Di-n-butyl phthalate | Ave | 1.380 | 1.370 | 0.0100 | 9.93 | 10.0 | -0.7 | 30.0 |
| Fluoranthene | Ave | 1.321 | 1.260 | 0.6000 | 9.54 | 10.0 | -4.6 | 30.0 |
| Benzidine | Ave | 0.5843 | 0.5955 | | 20.4 | 20.0 | 1.9 | 30.0 |
| Pyrene | Ave | 1.135 | 1.058 | 0.6000 | 9.33 | 10.0 | -6.7 | 30.0 |
| Butyl benzyl phthalate | Ave | 0.5142 | 0.5113 | 0.0100 | 9.94 | 10.0 | -0.6 | 30.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3952 | 0.3739 | 0.0100 | 18.9 | 20.0 | -5.4 | 30.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.6579 | 0.6437 | 0.0100 | 9.78 | 10.0 | -2.2 | 30.0 |
| Benzo[a]anthracene | Ave | 1.126 | 1.090 | 0.8000 | 9.68 | 10.0 | -3.2 | 30.0 |
| Chrysene | Ave | 1.076 | 1.070 | 0.7000 | 9.95 | 10.0 | -0.5 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-200091/11 | Calibration Date: 10/02/2015 14:28

Instrument ID: A4HP10 | Calib Start Date: 10/02/2015 10:48

GC Column: DB-624 ID: 0.18(mm) | Calib End Date: 10/02/2015 14:04

Lab File ID: 51002011.D | Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Ave | 1.371 | 1.402 | 0.0100 | 10.2 | 10.0 | 2.3 | 30.0 |
| Benzo[b]fluoranthene | Ave | 1.128 | 1.130 | 0.7000 | 10.0 | 10.0 | 0.2 | 30.0 |
| Benzo[k]fluoranthene | Ave | 1.274 | 1.168 | 0.7000 | 9.17 | 10.0 | -8.3 | 30.0 |
| Benzo[a]pyrene | Ave | 1.120 | 1.096 | 0.7000 | 9.78 | 10.0 | -2.2 | 30.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.233 | 1.234 | 0.5000 | 10.0 | 10.0 | 0.0 | 30.0 |
| Dibenz(a,h)anthracene | Ave | 1.024 | 1.069 | 0.4000 | 10.4 | 10.0 | 4.4 | 30.0 |
| Benzo[g,h,i]perylene | Ave | 1.092 | 1.107 | 0.5000 | 10.1 | 10.0 | 1.3 | 30.0 |
| 2-Fluorophenol (Surr) | Ave | 1.486 | 1.530 | | 10.3 | 10.0 | 3.0 | 30.0 |
| Phenol-d5 (Surr) | Ave | 1.814 | 1.897 | | 10.5 | 10.0 | 4.6 | 30.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.3982 | 0.4088 | | 10.3 | 10.0 | 2.6 | 30.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.297 | 1.284 | | 9.90 | 10.0 | -1.0 | 30.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1891 | 0.2114 | | 11.2 | 10.0 | 11.8 | 30.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7619 | 0.7747 | | 10.2 | 10.0 | 1.7 | 30.0 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 51111001.D DFTPP Injection Date: 11/11/2015

Instrument ID: A4HP10 DFTPP Injection Time: 09:12

Analysis Batch No.: 206256

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 40.9 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0)1 |
| 69 | Mass 69 relative abundance | 34.5 | |
| 70 | Less than 2.0% of mass 69 | 0.2 | (0.5)1 |
| 127 | 25.0 - 75.0% of mass 198 | 50.8 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.5 | |
| 275 | 10.0- 30.0% of mass 198 | 25.4 | |
| 365 | Greater than 0.75% of mass 198 | 3.2 | |
| 441 | Present, but less than mass 443 | 8.7 | |
| 442 | 40.0 - 110.0% of mass 198 | 95.2 | |
| 443 | 15.0 - 24.0% of mass 442 | 18.5 | (19.4)2 |

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCV 240-206256/2 | 51111002.D | 11/11/2015 | 09:34 |
| | MB 240-205786/23-A | 51111004.D | 11/11/2015 | 10:27 |
| | LCS 240-205786/24-A | 51111005.D | 11/11/2015 | 10:51 |
| F-VB-UST-1-14 | 240-57597-1 | 51111006.D | 11/11/2015 | 11:16 |
| F-VB-UST-2-14 | 240-57597-2 | 51111007.D | 11/11/2015 | 11:40 |
| F-VB-UST-3-14 | 240-57597-3 | 51111008.D | 11/11/2015 | 12:05 |
| F-VB-UST-4-14 | 240-57597-4 | 51111009.D | 11/11/2015 | 12:30 |
| F-VB-UST-5-14 | 240-57597-5 | 51111010.D | 11/11/2015 | 12:54 |
| F-VB-UST-6-14 | 240-57597-6 | 51111011.D | 11/11/2015 | 13:19 |
| F-VB-UST-7-14 | 240-57597-7 | 51111012.D | 11/11/2015 | 13:44 |
| F-VB-UST-8-14 | 240-57597-8 | 51111013.D | 11/11/2015 | 14:08 |
| F-VB-UST-9-14 | 240-57597-9 | 51111014.D | 11/11/2015 | 14:33 |
| F-VB-UST-10-14 | 240-57597-10 | 51111015.D | 11/11/2015 | 14:57 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206256/2  Calibration Date: 11/11/2015 09:34

Instrument ID: A4HP10  Calib Start Date: 10/02/2015 10:48

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/02/2015 14:04

Lab File ID: 51111002.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.6903 | 0.5705 | | 8.27 | 10.0 | -17.3 | 20.0 |
| N-Nitrosodimethylamine | Ave | 0.9253 | 0.7500 | | 8.11 | 10.0 | -18.9 | 20.0 |
| Pyridine | Ave | 1.684 | 1.387 | | 8.23 | 10.0 | -17.7 | 20.0 |
| Benzaldehyde | Ave | 1.111 | 0.6352 | 0.0100 | 11.4 | 20.0 | -42.8* | 20.0 |
| Phenol | Ave | 1.891 | 1.893 | 0.8000 | 10.0 | 10.0 | 0.0 | 20.0 |
| Aniline | Ave | 2.402 | 2.168 | | 9.03 | 10.0 | -9.7 | 20.0 |
| Bis(2-chloroethyl)ether | Ave | 1.506 | 1.420 | 0.7000 | 9.43 | 10.0 | -5.7 | 20.0 |
| 2-Chlorophenol | Ave | 1.543 | 1.438 | 0.8000 | 9.32 | 10.0 | -6.8 | 20.0 |
| n-Decane | Ave | 1.911 | 1.929 | | 10.1 | 10.0 | 0.9 | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.579 | 1.537 | | 9.73 | 10.0 | -2.7 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.615 | 1.575 | | 9.76 | 10.0 | -2.4 | 20.0 |
| Benzyl alcohol | Ave | 1.025 | 0.3563 | | 3.48 | 10.0 | -65.2* | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.531 | 1.486 | | 9.70 | 10.0 | -3.0 | 20.0 |
| 2-Methylphenol | Ave | 1.412 | 1.308 | 0.7000 | 9.27 | 10.0 | -7.3 | 20.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.121 | 2.250 | | 10.6 | 10.0 | 6.1 | 20.0 |
| Indene | Ave | 2.534 | 2.469 | | 19.5 | 20.0 | -2.6 | 20.0 |
| N-Nitrosodi-n-propylamine | Ave | 0.8785 | 0.8567 | 0.5000 | 9.75 | 10.0 | -2.5 | 20.0 |
| 3 & 4 Methylphenol | Ave | 1.469 | 1.380 | | 9.39 | 10.0 | -6.1 | 20.0 |
| Acetophenone | Ave | 1.988 | 1.975 | 0.0100 | 9.93 | 10.0 | -0.7 | 20.0 |
| Hexachloroethane | Ave | 0.6359 | 0.6671 | 0.3000 | 10.5 | 10.0 | 4.9 | 20.0 |
| Nitrobenzene | Ave | 0.3694 | 0.3589 | 0.2000 | 9.72 | 10.0 | -2.8 | 20.0 |
| Isophorone | Ave | 0.6852 | 0.6730 | 0.4000 | 9.82 | 10.0 | -1.8 | 20.0 |
| 2-Nitrophenol | Ave | 0.1669 | 0.1844 | 0.1000 | 11.0 | 10.0 | 10.5 | 20.0 |
| 2,4-Dimethylphenol | Ave | 0.3580 | 0.3249 | 0.2000 | 9.07 | 10.0 | -9.3 | 20.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4236 | 0.4023 | 0.3000 | 9.50 | 10.0 | -5.0 | 20.0 |
| Benzoic acid | Lin1 | | 0.1787 | | 17.0 | 20.0 | -15.1 | 20.0 |
| 2,4-Dichlorophenol | Ave | 0.2940 | 0.2941 | 0.2000 | 10.0 | 10.0 | 0.0 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.3022 | 0.3090 | | 10.2 | 10.0 | 2.2 | 20.0 |
| Naphthalene | Ave | 1.084 | 1.060 | 0.7000 | 9.78 | 10.0 | -2.2 | 20.0 |
| 4-Chloroaniline | Ave | 0.4756 | 0.4439 | 0.0100 | 9.33 | 10.0 | -6.7 | 20.0 |
| 2,6-Dichlorophenol | Ave | 0.2949 | 0.2816 | | 9.55 | 10.0 | -4.5 | 20.0 |
| Hexachlorobutadiene | Ave | 0.1868 | 0.1936 | 0.0100 | 10.4 | 10.0 | 3.6 | 20.0 |
| Caprolactam | Ave | 0.1164 | 0.1013 | 0.0100 | 17.4 | 20.0 | -12.9 | 20.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3033 | 0.3014 | 0.2000 | 9.94 | 10.0 | -0.6 | 20.0 |
| 2-Methylnaphthalene | Ave | 0.7006 | 0.6945 | 0.4000 | 9.91 | 10.0 | -0.9 | 20.0 |
| 1-Methylnaphthalene | Ave | 0.6254 | 0.6095 | | 9.75 | 10.0 | -2.5 | 20.0 |
| Hexachlorocyclopentadiene | Ave | 0.2689 | 0.1923 | 0.0500 | 7.15 | 10.0 | -28.5* | 20.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4998 | 0.5133 | 0.0100 | 10.3 | 10.0 | 2.7 | 20.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3619 | 0.3550 | 0.2000 | 9.81 | 10.0 | -1.9 | 20.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3794 | 0.3767 | 0.2000 | 9.93 | 10.0 | -0.7 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206256/2  Calibration Date: 11/11/2015 09:34

Instrument ID: A4HP10  Calib Start Date: 10/02/2015 10:48

GC Column: DB-624  ID: 0.18(mm)  Calib End Date: 10/02/2015 14:04

Lab File ID: 51111002.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.409 | 1.370 | 0.0100 | 9.72 | 10.0 | -2.8 | 20.0 |
| 2-Chloronaphthalene | Ave | 1.182 | 1.123 | 0.8000 | 9.50 | 10.0 | -5.0 | 20.0 |
| 2-Nitroaniline | Ave | 0.3503 | 0.3581 | 0.0100 | 10.2 | 10.0 | 2.2 | 20.0 |
| Dimethyl phthalate | Ave | 1.285 | 1.285 | 0.0100 | 10.0 | 10.0 | 0.0 | 20.0 |
| 1,3-Dinitrobenzene | Lin1 | | 0.2007 | | 9.65 | 10.0 | -3.5 | 20.0 |
| 2,6-Dinitrotoluene | Ave | 0.2818 | 0.2901 | | 10.3 | 10.0 | 3.0 | 20.0 |
| Acenaphthylene | Ave | 1.763 | 1.739 | 0.9000 | 9.87 | 10.0 | -1.3 | 20.0 |
| 3-Nitroaniline | Ave | 0.3239 | 0.3137 | 0.0100 | 9.68 | 10.0 | -3.2 | 20.0 |
| Acenaphthene | Ave | 1.075 | 1.050 | 0.9000 | 9.77 | 10.0 | -2.3 | 20.0 |
| 2,4-Dinitrophenol | Lin1 | | 0.1133 | 0.0100 | 16.5 | 20.0 | -17.5 | 20.0 |
| 4-Nitrophenol | Ave | 0.1702 | 0.1831 | | 21.5 | 20.0 | 7.6 | 20.0 |
| 2,4-Dinitrotoluene | Ave | 0.3666 | 0.3849 | 0.2000 | 10.5 | 10.0 | 5.0 | 20.0 |
| Dibenzofuran | Ave | 1.623 | 1.575 | 0.8000 | 9.70 | 10.0 | -3.0 | 20.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.3179 | 0.2913 | 0.0100 | 9.16 | 10.0 | -8.4 | 20.0 |
| Diethyl phthalate | Ave | 1.172 | 1.234 | 0.0100 | 10.5 | 10.0 | 5.3 | 20.0 |
| Hexadecane | Ave | 0.5546 | 0.5740 | | 10.3 | 10.0 | 3.5 | 20.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.6110 | 0.5883 | 0.4000 | 9.63 | 10.0 | -3.7 | 20.0 |
| 4-Nitroaniline | Ave | 0.3149 | 0.2932 | 0.0100 | 9.31 | 10.0 | -6.9 | 20.0 |
| Fluorene | Ave | 1.322 | 1.299 | 0.9000 | 9.82 | 10.0 | -1.8 | 20.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1067 | 0.0100 | 16.0 | 20.0 | -20.0 | 20.0 |
| N-Nitrosodiphenylamine | Ave | 0.5710 | 0.5894 | 0.0100 | 20.6 | 20.0 | 3.2 | 20.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 0.7994 | 0.7848 | | 9.81 | 10.0 | -1.8 | 20.0 |
| Azobenzene | Ave | 0.7997 | 0.7841 | | 9.80 | 10.0 | -2.0 | 20.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.2298 | 0.2307 | 0.1000 | 10.0 | 10.0 | 0.4 | 20.0 |
| Hexachlorobenzene | Ave | 0.2582 | 0.2467 | 0.1000 | 9.55 | 10.0 | -4.5 | 20.0 |
| Atrazine | Ave | 0.2140 | 0.2111 | 0.0100 | 19.7 | 20.0 | -1.4 | 20.0 |
| n-Octadecane | Qua | | 0.9708 | | 13.8 | 10.0 | 37.9* | 20.0 |
| Pentachlorophenol | Ave | 0.1432 | 0.1286 | 0.0500 | 18.0 | 20.0 | -10.2 | 20.0 |
| Phenanthrene | Ave | 1.177 | 1.137 | 0.7000 | 9.66 | 10.0 | -3.4 | 20.0 |
| Anthracene | Ave | 1.190 | 1.146 | 0.7000 | 9.63 | 10.0 | -3.7 | 20.0 |
| Carbazole | Ave | 1.127 | 1.044 | 0.0100 | 9.26 | 10.0 | -7.4 | 20.0 |
| Di-n-butyl phthalate | Ave | 1.380 | 1.388 | 0.0100 | 10.1 | 10.0 | 0.6 | 20.0 |
| Fluoranthene | Ave | 1.321 | 1.235 | 0.6000 | 9.35 | 10.0 | -6.5 | 20.0 |
| Benzidine | Ave | 0.5843 | 0.3860 | | 13.2 | 20.0 | -33.9* | 20.0 |
| Pyrene | Ave | 1.135 | 1.113 | 0.6000 | 9.81 | 10.0 | -1.9 | 20.0 |
| Butyl benzyl phthalate | Ave | 0.5142 | 0.5301 | 0.0100 | 10.3 | 10.0 | 3.1 | 20.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.6579 | 0.7290 | 0.0100 | 11.1 | 10.0 | 10.8 | 20.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3952 | 0.3826 | 0.0100 | 19.4 | 20.0 | -3.2 | 20.0 |
| Benzo[a]anthracene | Ave | 1.126 | 1.060 | 0.8000 | 9.41 | 10.0 | -5.9 | 20.0 |
| Chrysene | Ave | 1.076 | 0.9787 | 0.7000 | 9.10 | 10.0 | -9.0 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206256/2 | Calibration Date: 11/11/2015 09:34

Instrument ID: A4HP10 | Calib Start Date: 10/02/2015 10:48

GC Column: DB-624 ID: 0.18(mm) | Calib End Date: 10/02/2015 14:04

Lab File ID: 51111002.D | Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Ave | 1.371 | 1.497 | 0.0100 | 10.9 | 10.0 | 9.2 | 20.0 |
| Benzo[b]fluoranthene | Ave | 1.128 | 1.167 | 0.7000 | 10.3 | 10.0 | 3.5 | 20.0 |
| Benzo[k]fluoranthene | Ave | 1.274 | 1.217 | 0.7000 | 9.55 | 10.0 | -4.5 | 20.0 |
| Benzo[a]pyrene | Ave | 1.120 | 1.130 | 0.7000 | 10.1 | 10.0 | 0.9 | 20.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.233 | 1.265 | 0.5000 | 10.3 | 10.0 | 2.6 | 20.0 |
| Dibenz(a,h)anthracene | Ave | 1.024 | 1.075 | 0.4000 | 10.5 | 10.0 | 5.0 | 20.0 |
| Benzo[g,h,i]perylene | Ave | 1.092 | 1.055 | 0.5000 | 9.66 | 10.0 | -3.4 | 20.0 |
| 2-Fluorophenol (Surr) | Ave | 1.486 | 1.313 | | 8.83 | 10.0 | -11.7 | 20.0 |
| Phenol-d5 (Surr) | Ave | 1.814 | 1.632 | | 9.00 | 10.0 | -10.0 | 20.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.3982 | 0.3544 | | 8.90 | 10.0 | -11.0 | 20.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.297 | 1.228 | | 9.47 | 10.0 | -5.3 | 20.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1891 | 0.1700 | | 8.99 | 10.0 | -10.1 | 20.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7619 | 0.7712 | | 10.1 | 10.0 | 1.2 | 20.0 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab File ID: 51030001.D DFTPP Injection Date: 10/30/2015

Instrument ID: A4HP9 DFTPP Injection Time: 12:57

Analysis Batch No.: 204569

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 46.3 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0)1 |
| 69 | Mass 69 relative abundance | 51.4 | |
| 70 | Less than 2.0% of mass 69 | 0.2 | (0.4)1 |
| 127 | 25.0 - 75.0% of mass 198 | 59.4 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.3 | |
| 275 | 10.0- 30.0% of mass 198 | 21.2 | |
| 365 | Greater than 0.75% of mass 198 | 3.2 | |
| 441 | Present, but less than mass 443 | 0.3 | |
| 442 | 40.0 - 110.0% of mass 198 | 85.4 | |
| 443 | 15.0 - 24.0% of mass 442 | 16.8 | (19.7)2 |

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | STD5 240-204569/2 | 51030002.D | 10/30/2015 | 13:19 |
| | STD4 240-204569/3 | 51030003.D | 10/30/2015 | 13:44 |
| | STD3 240-204569/4 | 51030004.D | 10/30/2015 | 14:10 |
| | STD2 240-204569/5 | 51030005.D | 10/30/2015 | 14:35 |
| | STD1 240-204569/6 | 51030006.D | 10/30/2015 | 15:01 |
| | STD9 240-204569/7 | 51030007.D | 10/30/2015 | 15:27 |
| | STD8 240-204569/8 | 51030008.D | 10/30/2015 | 15:52 |
| | STD7 240-204569/9 | 51030009.D | 10/30/2015 | 16:17 |
| | STD6 240-204569/10 | 51030010.D | 10/30/2015 | 16:42 |
| | ICV 240-204569/11 | 51030011.D | 10/30/2015 | 17:08 |

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19  Calibration End Date: 10/30/2015 16:42  Calibration ID: 31170

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-204569/6 | 51030006.D |
| Level 2 | STD2 240-204569/5 | 51030005.D |
| Level 3 | STD3 240-204569/4 | 51030004.D |
| Level 4 | STD4 240-204569/3 | 51030003.D |
| Level 5 | STD5 240-204569/2 | 51030002.D |
| Level 6 | STD6 240-204569/10 | 51030010.D |
| Level 7 | STD7 240-204569/9 | 51030009.D |
| Level 8 | STD8 240-204569/8 | 51030008.D |
| Level 9 | STD9 240-204569/7 | 51030007.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 1,4-Dioxane | 0.7879 | 0.6702<br>0.7724 | 0.7963<br>0.7912 | 0.7519<br>0.7577 | 0.7576 | Ave | | 0.7606 | | | | 5.3 | | 20.0 | | | |
| N-Nitrosodimethylamine | 1.1590 | 1.0469<br>1.1257 | 0.9916<br>1.1673 | 1.0993<br>1.1059 | 1.0809 | Ave | | 1.0971 | | | | 5.3 | | 20.0 | | | |
| Pyridine | 2.0907 | 2.0205<br>1.9910 | 1.8876<br>2.0977 | 2.0118<br>1.9855 | 2.0565 | Ave | | 2.0177 | | | | 3.3 | | 20.0 | | | |
| Benzaldehyde | 1.3725 | 1.5727<br>1.2076 | 1.5657<br>1.2344 | 1.5133<br>+++++ | 1.4739 | Ave | | 1.4200 | | | 0.0100 | 10.7 | | 20.0 | | | |
| Phenol | 2.1853 | 2.1727<br>2.0136 | 2.0934<br>2.0804 | 2.1195<br>1.9722 | 2.1525 | Ave | | 2.0987 | | | 0.8000 | 3.6 | | 20.0 | | | |
| Aniline | 2.7280 | 2.6247<br>2.5385 | 2.6136<br>2.6466 | 2.6262<br>2.5459 | 2.6111 | Ave | | 2.6168 | | | | 2.3 | | 20.0 | | | |
| Bis(2-chloroethyl)ether | 1.8558 | 1.9142<br>1.6942 | 1.8429<br>1.7572 | 1.8447<br>1.6890 | 1.8270 | Ave | | 1.8031 | | | 0.7000 | 4.5 | | 20.0 | | | |
| 2-Chlorophenol | 1.5989 | 1.4425<br>1.5006 | 1.5455<br>1.5432 | 1.5087<br>1.4819 | 1.5120 | Ave | | 1.5167 | | | 0.8000 | 3.1 | | 20.0 | | | |
| n-Decane | 1.7829 | 1.8718<br>1.6389 | 1.8936<br>1.6552 | 1.8312<br>1.5772 | 1.8099 | Ave | | 1.7576 | | | | 6.7 | | 20.0 | | | |
| 1,3-Dichlorobenzene | 1.6084 | 1.5596<br>1.5011 | 1.6679<br>1.5578 | 1.6175<br>1.4910 | 1.5792 | Ave | | 1.5728 | | | | 3.8 | | 20.0 | | | |
| 1,4-Dichlorobenzene | 1.6334 | 1.6944<br>1.5231 | 1.6061<br>1.5860 | 1.6427<br>1.5227 | 1.6060 | Ave | | 1.6018 | | | | 3.7 | | 20.0 | | | |
| Benzyl alcohol | 1.1163 | 0.8791<br>1.0801 | 0.9185<br>1.1234 | 0.9897<br>1.0846 | 1.0413 | Ave | | 1.0291 | | | | 8.9 | | 20.0 | | | |
| 1,2-Dichlorobenzene | 1.5481 | 1.6647<br>1.4507 | 1.5548<br>1.5234 | 1.5786<br>1.4511 | 1.5515 | Ave | | 1.5404 | | | | 4.5 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2-Methylphenol | 1.6435 | 1.5529<br>1.5278 | 1.5884<br>1.6065 | 1.5262<br>1.5342 | 1.5781 | Ave | | 1.5697 | | | 0.7000 | 2.7 | | 20.0 | | | |
| bis (2-chloroisopropyl) ether | 2.0663 | 2.1290<br>1.9114 | 2.1504<br>1.9756 | 2.0501<br>1.8678 | 2.0388 | Ave | | 2.0237 | | | | 4.9 | | 20.0 | | | |
| Indene | 2.7276 | 2.8401<br>2.4658 | 2.8969<br>2.5378 | 2.8654<br>2.3993 | 2.7687 | Ave | | 2.6877 | | | | 7.2 | | 20.0 | | | |
| 3 & 4 Methylphenol | 1.6857 | 1.6121<br>1.5969 | 1.5331<br>1.6527 | 1.6625<br>1.5643 | 1.6490 | Ave | | 1.6195 | | | | 3.2 | | 20.0 | | | |
| N-Nitrosodi-n-propylamine | 1.1989 | 1.1171<br>1.1083 | 1.1602<br>1.1430 | 1.1470<br>1.0752 | 1.1583 | Ave | | 1.1385 | | | 0.5000 | 3.3 | | 20.0 | | | |
| Acetophenone | 2.3832 | 2.4667<br>2.2595 | 2.3410<br>2.3395 | 2.3767<br>2.2604 | 2.3981 | Ave | | 2.3531 | | | 0.0100 | 3.0 | | 20.0 | | | |
| Hexachloroethane | 0.6860 | 0.6898<br>0.6531 | 0.6582<br>0.6793 | 0.6592<br>0.6596 | 0.6614 | Ave | | 0.6683 | | | 0.3000 | 2.1 | | 20.0 | | | |
| Nitrobenzene | 0.4587 | 0.4485<br>0.4327 | 0.4414<br>0.4340 | 0.4609<br>0.4248 | 0.4662 | Ave | | 0.4459 | | | 0.2000 | 3.4 | | 20.0 | | | |
| Isophorone | 0.7828 | 0.7464<br>0.7590 | 0.7806<br>0.7476 | 0.7726<br>0.7157 | 0.7963 | Ave | | 0.7626 | | | 0.4000 | 3.4 | | 20.0 | | | |
| 2-Nitrophenol | 0.1828 | 0.1373<br>0.1790 | 0.1472<br>0.1753 | 0.1582<br>0.1700 | 0.1781 | Ave | | 0.1660 | | | 0.1000 | 10.0 | | 20.0 | | | |
| 2,4-Dimethylphenol | 0.3869 | 0.3987<br>0.3723 | 0.4012<br>0.3647 | 0.3813<br>0.3471 | 0.4020 | Ave | | 0.3818 | | | 0.2000 | 5.1 | | 20.0 | | | |
| Benzoic acid | 0.2127 | +++++<br>0.2174 | 0.0852<br>0.2213 | 0.1129<br>0.2114 | 0.1510 | Lin1 | -0.359 | 0.2234 | | | | | | | 0.9960 | | 0.9900 |
| Bis(2-chloroethoxy)methane | 0.4632 | 0.4830<br>0.4267 | 0.4845<br>0.4231 | 0.4695<br>0.4030 | 0.4779 | Ave | | 0.4539 | | | 0.3000 | 7.0 | | 20.0 | | | |
| 2,4-Dichlorophenol | 0.2586 | 0.2322<br>0.2510 | 0.2551<br>0.2491 | 0.2461<br>0.2424 | 0.2571 | Ave | | 0.2489 | | | 0.2000 | 3.5 | | 20.0 | | | |
| 1,2,4-Trichlorobenzene | 0.2904 | 0.3057<br>0.2778 | 0.3084<br>0.2753 | 0.2978<br>0.2688 | 0.3048 | Ave | | 0.2911 | | | | 5.3 | | 20.0 | | | |
| Naphthalene | 1.3148<br>1.0745 | 1.1643<br>1.0146 | 1.1361<br>0.9888 | 1.0999<br>0.9493 | 1.0983 | Ave | | 1.0934 | | | 0.7000 | 9.9 | | 20.0 | | | |
| 4-Chloroaniline | 0.4907 | 0.4694<br>0.4709 | 0.4675<br>0.4648 | 0.4499<br>0.4471 | 0.4606 | Ave | | 0.4651 | | | 0.0100 | 2.9 | | 20.0 | | | |
| 2,6-Dichlorophenol | 0.2619 | 0.2609<br>0.2504 | 0.2649<br>0.2440 | 0.2474<br>0.2401 | 0.2580 | Ave | | 0.2535 | | | | 3.6 | | 20.0 | | | |
| Hexachlorobutadiene | 0.1543 | 0.1757<br>0.1478 | 0.1596<br>0.1459 | 0.1568<br>0.1417 | 0.1563 | Ave | | 0.1548 | | | 0.0100 | 6.8 | | 20.0 | | | |
| Caprolactam | 0.1199 | 0.0914<br>0.1182 | 0.1059<br>0.1173 | 0.1129<br>0.1147 | 0.1141 | Ave | | 0.1118 | | | 0.0100 | 8.3 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9 GC Column: RXI-5SILMS ID: 0.45(mm) Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19 Calibration End Date: 10/30/2015 16:42 Calibration ID: 31170

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 4-Chloro-3-methylphenol | 0.3405 | 0.2915<br>0.3291 | 0.2982<br>0.3289 | 0.3051<br>0.3200 | 0.3280 | Ave | | 0.3177 | | | 0.2000 | 5.5 | | 20.0 | | | |
| 2-Methylnaphthalene | 0.6714<br>0.6849 | 0.7552<br>0.6463 | 0.7151<br>0.6273 | 0.6901<br>0.5996 | 0.7011 | Ave | | 0.6768 | | | 0.4000 | 7.0 | | 20.0 | | | |
| 1-Methylnaphthalene | 0.6508<br>0.6069 | 0.6257<br>0.5675 | 0.6232<br>0.5547 | 0.6191<br>0.5412 | 0.6146 | Ave | | 0.6004 | | | | 6.2 | | 20.0 | | | |
| Hexachlorocyclopentadiene | 0.2836 | 0.2033<br>0.2937 | 0.2109<br>0.2923 | 0.2357<br>0.2863 | 0.2621 | Ave | | 0.2585 | | | 0.0500 | 14.3 | | 20.0 | | | |
| 1,2,4,5-Tetrachlorobenzene | 0.4829 | 0.5543<br>0.4681 | 0.5241<br>0.4467 | 0.5277<br>0.4361 | 0.5085 | Ave | | 0.4935 | | | 0.0100 | 8.5 | | 20.0 | | | |
| 2,4,6-Trichlorophenol | 0.3520 | 0.3202<br>0.3511 | 0.3243<br>0.3405 | 0.3380<br>0.3354 | 0.3397 | Ave | | 0.3377 | | | 0.2000 | 3.3 | | 20.0 | | | |
| 2,4,5-Trichlorophenol | 0.3611 | 0.3258<br>0.3643 | 0.3548<br>0.3514 | 0.3345<br>0.3507 | 0.3665 | Ave | | 0.3511 | | | 0.2000 | 4.1 | | 20.0 | | | |
| 1,1'-Biphenyl | 1.7382<br>1.5054 | 1.7110<br>1.4492 | 1.5822<br>1.3989 | 1.5786<br>1.3700 | 1.5456 | Ave | | 1.5421 | | | 0.0100 | 8.3 | | 20.0 | | | |
| 2-Chloronaphthalene | 1.3119<br>1.1298 | 1.2573<br>1.0949 | 1.1833<br>1.0732 | 1.1920<br>1.0363 | 1.1795 | Ave | | 1.1620 | | | 0.8000 | 7.6 | | 20.0 | | | |
| 2-Nitroaniline | 0.4424 | 0.3228<br>0.4442 | 0.3489<br>0.4281 | 0.3983<br>0.4198 | 0.4199 | Ave | | 0.4031 | | | 0.0100 | 11.0 | | 20.0 | | | |
| Dimethyl phthalate | 1.2250 | 1.3046<br>1.1901 | 1.2480<br>1.1306 | 1.2490<br>1.1063 | 1.2488 | Ave | | 1.2128 | | | 0.0100 | 5.5 | | 20.0 | | | |
| 1,3-Dinitrobenzene | 0.2148 | 0.1418<br>0.2208 | 0.1424<br>0.2112 | 0.1739<br>0.2124 | 0.1993 | Ave | | 0.1896 | | | | 17.2 | | 20.0 | | | |
| 2,6-Dinitrotoluene | 0.3015 | 0.2515<br>0.2974 | 0.2744<br>0.2833 | 0.2734<br>0.2819 | 0.2878 | Ave | | 0.2814 | | | | 5.6 | | 20.0 | | | |
| Acenaphthylene | 2.1239<br>1.9191 | 1.8953<br>1.8490 | 1.9145<br>1.7759 | 1.9143<br>1.7344 | 1.9293 | Ave | | 1.8951 | | | 0.9000 | 5.8 | | 20.0 | | | |
| 3-Nitroaniline | 0.3684 | 0.2911<br>0.3658 | 0.3323<br>0.3564 | 0.3421<br>0.3575 | 0.3522 | Ave | | 0.3457 | | | 0.0100 | 7.2 | | 20.0 | | | |
| 2,4-Dinitrophenol | 0.1226 | +++++<br>0.1490 | 0.0184<br>0.1495 | 0.0295<br>0.1498 | 0.0689 | Lin | -0.576 | 0.1620 | | | 0.0100 | | | | 0.9940 | | 0.9900 |
| Acenaphthene | 1.4818<br>1.1919 | 1.3172<br>1.1478 | 1.2378<br>1.0901 | 1.2469<br>1.0790 | 1.2360 | Ave | | 1.2254 | | | 0.9000 | 10.1 | | 20.0 | | | |
| 4-Nitrophenol | 0.2363 | 0.1350<br>0.2358 | 0.1725<br>0.2308 | 0.1994<br>0.2253 | 0.2085 | Ave | | 0.2054 | | | | 17.5 | | 20.0 | | | |
| 2,4-Dinitrotoluene | 0.3819 | 0.2653<br>0.3890 | 0.3195<br>0.3921 | 0.3356<br>0.3853 | 0.3775 | Ave | | 0.3558 | | | 0.2000 | 12.7 | | 20.0 | | | |
| Dibenzofuran | 1.9633<br>1.5887 | 1.8065<br>1.5423 | 1.6844<br>1.4828 | 1.6535<br>1.4457 | 1.6428 | Ave | | 1.6455 | | | 0.8000 | 9.8 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19  Calibration End Date: 10/30/2015 16:42  Calibration ID: 31170

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| 2,3,4,6-Tetrachlorophenol | 0.2882 | 0.2083<br>0.2752 | 0.2408<br>0.2694 | 0.2697<br>0.2698 | 0.2803 | Ave | | 0.2627 | | | 0.0100 | 9.9 | | 20.0 | | | |
| Hexadecane | 0.8423 | 0.9528<br>0.7419 | 0.9952<br>0.6898 | 0.9799<br>+++++ | 0.9285 | Ave | | 0.8758 | | | | 13.8 | | 20.0 | | | |
| Diethyl phthalate | 1.1435 | 1.3219<br>1.1140 | 1.3222<br>1.0714 | 1.3321<br>1.0617 | 1.2828 | Ave | | 1.2062 | | | 0.0100 | 9.9 | | 20.0 | | | |
| 4-Chlorophenyl phenyl ether | 0.5893 | 0.6305<br>0.5829 | 0.6266<br>0.5548 | 0.6099<br>0.5380 | 0.5949 | Ave | | 0.5909 | | | 0.4000 | 5.5 | | 20.0 | | | |
| 4-Nitroaniline | 0.3862 | 0.2821<br>0.3889 | 0.3158<br>0.3676 | 0.3459<br>0.3461 | 0.3410 | Ave | | 0.3467 | | | 0.0100 | 10.3 | | 20.0 | | | |
| Fluorene | 1.4143<br>1.3000 | 1.3977<br>1.2358 | 1.3185<br>1.1690 | 1.3302<br>1.1121 | 1.3108 | Ave | | 1.2876 | | | 0.9000 | 7.7 | | 20.0 | | | |
| 4,6-Dinitro-2-methylphenol | 0.1273 | +++++<br>0.1312 | 0.0456<br>0.1282 | 0.0664<br>0.1258 | 0.0946 | Lin1 | -0.215 | 0.1327 | | | 0.0100 | | | | 0.9970 | | 0.9900 |
| N-Nitrosodiphenylamine | 0.5686 | 0.5541<br>0.4887 | 0.5453<br>0.4764 | 0.5445<br>0.4597 | 0.5726 | Ave | | 0.5262 | | | 0.0100 | 8.4 | | 20.0 | | | |
| 1,2-Diphenylhydrazine(as Azobenzene) | 1.1353 | 1.0022<br>0.9923 | 0.9840<br>0.9576 | 1.0232<br>0.9032 | 1.1113 | Ave | | 1.0141 | | | | 7.6 | | 20.0 | | | |
| Azobenzene | 1.1352 | 1.0064<br>0.9924 | 0.9829<br>0.9577 | 1.0228<br>0.9034 | 1.1118 | Ave | | 1.0141 | | | | 7.6 | | 20.0 | | | |
| 4-Bromophenyl phenyl ether | 0.2076 | 0.1841<br>0.1792 | 0.1926<br>0.1792 | 0.1849<br>0.1705 | 0.1914 | Ave | | 0.1862 | | | 0.1000 | 6.0 | | 20.0 | | | |
| Hexachlorobenzene | 0.2326<br>0.2013 | 0.2025<br>0.1742 | 0.1926<br>0.1716 | 0.1883<br>0.1642 | 0.1879 | Ave | | 0.1906 | | | 0.1000 | 10.8 | | 20.0 | | | |
| Atrazine | 0.2016 | 0.1731<br>0.1708 | 0.1758<br>0.1648 | 0.1724<br>0.1578 | 0.1894 | Ave | | 0.1757 | | | 0.0100 | 7.9 | | 20.0 | | | |
| n-Octadecane | 0.5786 | 0.4969<br>0.4950 | 0.5220<br>0.4662 | 0.5373<br>0.4319 | 0.5750 | Ave | | 0.5128 | | | | 9.9 | | 20.0 | | | |
| Pentachlorophenol | 0.1310 | 0.0630<br>0.1277 | 0.0723<br>0.1247 | 0.0870<br>0.1199 | 0.1058 | Lin1 | -0.089 | 0.1258 | | | 0.0500 | | | | 0.9970 | | 0.9900 |
| Phenanthrene | 1.3770<br>1.1812 | 1.2617<br>1.0880 | 1.2072<br>1.0839 | 1.1682<br>1.0482 | 1.1923 | Ave | | 1.1786 | | | 0.7000 | 8.6 | | 20.0 | | | |
| Anthracene | 1.1316<br>1.2010 | 1.1640<br>1.1155 | 1.1522<br>1.1046 | 1.1498<br>1.0751 | 1.1966 | Ave | | 1.1434 | | | 0.7000 | 3.6 | | 20.0 | | | |
| Carbazole | 1.1181 | 0.9855<br>1.0121 | 1.0010<br>0.9736 | 0.9681<br>0.9190 | 0.9888 | Ave | | 0.9958 | | | 0.0100 | 5.7 | | 20.0 | | | |
| Di-n-butyl phthalate | 1.4278 | 1.0543<br>1.2449 | 1.0537<br>1.1894 | 1.0836<br>1.1345 | 1.2603 | Ave | | 1.1811 | | | 0.0100 | 10.9 | | 20.0 | | | |
| Fluoranthene | 1.1437<br>1.2640 | 1.1063<br>1.1301 | 1.1177<br>1.0882 | 1.0757<br>1.0483 | 1.1538 | Ave | | 1.1253 | | | 0.6000 | 5.5 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 204569

SDG No.:

Instrument ID: A4HP9  GC Column: RXI-5SILMS  ID: 0.45(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 10/30/2015 13:19  Calibration End Date: 10/30/2015 16:42  Calibration ID: 31170

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2<br>LVL 7 | LVL 3<br>LVL 8 | LVL 4<br>LVL 9 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| Benzidine | 0.6951 | 0.3632<br>0.6772 | 0.4149<br>0.6916 | 0.4087<br>0.6334 | 0.5454 | Lin1 | -0.483 | 0.6732 | | | | | | | 0.9940 | | 0.9900 |
| Pyrene | 1.2053<br>1.1794 | 1.2231<br>1.1038 | 1.2032<br>1.1259 | 1.1539<br>1.0853 | 1.2062 | Ave | | 1.1651 | | | 0.6000 | 4.3 | | 20.0 | | | |
| Butyl benzyl phthalate | 0.5806 | 0.4429<br>0.5620 | 0.4886<br>0.5776 | 0.4908<br>0.5538 | 0.5224 | Ave | | 0.5273 | | | 0.0100 | 9.4 | | 20.0 | | | |
| Bis(2-ethylhexyl) phthalate | 0.7846 | 0.5860<br>0.7590 | 0.6528<br>0.7817 | 0.6396<br>0.7552 | 0.7101 | Ave | | 0.7086 | | | 0.0100 | 10.5 | | 20.0 | | | |
| 3,3'-Dichlorobenzidine | 0.3701 | 0.2827<br>0.3612 | 0.3127<br>0.3755 | 0.3218<br>0.3615 | 0.3555 | Ave | | 0.3426 | | | 0.0100 | 9.6 | | 20.0 | | | |
| Benzo[a]anthracene | 1.1876<br>1.1606 | 1.1259<br>1.0983 | 1.0942<br>1.1316 | 1.0394<br>1.1004 | 1.0862 | Ave | | 1.1138 | | | 0.8000 | 3.9 | | 20.0 | | | |
| Chrysene | 1.2471<br>1.1079 | 1.2562<br>1.0636 | 1.1353<br>1.0780 | 1.0870<br>1.0408 | 1.1277 | Ave | | 1.1271 | | | 0.7000 | 6.8 | | 20.0 | | | |
| Di-n-octyl phthalate | 1.5859 | 0.8921<br>1.6214 | 1.0660<br>1.6195 | 1.1495<br>1.5780 | 1.2347 | Lin1 | -0.537 | 1.6089 | | | 0.0100 | | | | 0.9960 | | 0.9900 |
| Benzo[b]fluoranthene | 0.6162<br>1.2204 | 0.9128<br>1.1527 | 0.9577<br>1.1512 | 1.0001<br>1.2028 | 1.0888 | Ave | | 1.0336 | | | 0.7000 | 18.4 | | 20.0 | | | |
| Benzo[k]fluoranthene | 1.3131<br>1.3193 | 1.3524<br>1.3266 | 1.2758<br>1.3571 | 1.3772<br>1.2656 | 1.4179 | Ave | | 1.3339 | | | 0.7000 | 3.6 | | 20.0 | | | |
| Benzo[a]pyrene | 0.8263<br>1.2313 | 1.0262<br>1.1934 | 1.0302<br>1.1949 | 1.0741<br>1.1922 | 1.1168 | Ave | | 1.0984 | | | 0.7000 | 11.6 | | 20.0 | | | |
| Indeno[1,2,3-cd]pyrene | 0.7163<br>1.2561 | 0.9705<br>1.2084 | 0.9906<br>1.2361 | 1.1154<br>1.2234 | 1.2408 | Ave | | 1.1064 | | | 0.5000 | 16.5 | | 20.0 | | | |
| Dibenz(a,h)anthracene | 0.5695<br>1.0470 | 0.7363<br>0.9974 | 0.7350<br>1.0228 | 0.8353<br>0.9987 | 0.9400 | Ave | | 0.8758 | | | 0.4000 | 18.9 | | 20.0 | | | |
| Benzo[g,h,i]perylene | 0.8647<br>1.0804 | 0.9477<br>1.0159 | 0.9127<br>1.0424 | 0.9792<br>1.0274 | 1.0726 | Ave | | 0.9937 | | | 0.5000 | 7.4 | | 20.0 | | | |
| 2-Fluorophenol (Surr) | 1.5911 | 1.4167<br>1.5027 | 1.4944<br>1.5649 | 1.5041<br>1.4888 | 1.5108 | Ave | | 1.5092 | | | | 3.5 | | 20.0 | | | |
| Phenol-d5 (Surr) | 2.0299 | 1.9547<br>1.8921 | 2.0082<br>1.9505 | 2.0039<br>1.8641 | 1.9846 | Ave | | 1.9610 | | | | 3.0 | | 20.0 | | | |
| Nitrobenzene-d5 (Surr) | 0.5016<br>0.4763 | 0.4486<br>0.4608 | 0.4786<br>0.4646 | 0.4647<br>0.4426 | 0.4825 | Ave | | 0.4689 | | | | 3.9 | | 20.0 | | | |
| 2-Fluorobiphenyl (Surr) | 1.4656<br>1.2301 | 1.3221<br>1.1942 | 1.2848<br>1.1497 | 1.3000<br>1.1252 | 1.2728 | Ave | | 1.2605 | | | | 8.1 | | 20.0 | | | |
| 2,4,6-Tribromophenol (Surr) | 0.1475 | 0.1223<br>0.1518 | 0.1207<br>0.1515 | 0.1348<br>0.1508 | 0.1415 | Ave | | 0.1401 | | | | 9.2 | | 20.0 | | | |
| Terphenyl-d14 (Surr) | 0.8153<br>0.7449 | 0.7288<br>0.7126 | 0.7158<br>0.7256 | 0.7120<br>0.7014 | 0.7254 | Ave | | 0.7313 | | | | 4.6 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204569/11 | Calibration Date: 10/30/2015 17:08

Instrument ID: A4HP9 | Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS ID: 0.45(mm) | Calib End Date: 10/30/2015 16:42

Lab File ID: 51030011.D | Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,4-Dioxane | Ave | 0.7606 | 0.7555 | | 9.93 | 10.0 | -0.7 | 30.0 |
| N-Nitrosodimethylamine | Ave | 1.097 | 0.9653 | | 8.80 | 10.0 | -12.0 | 30.0 |
| Pyridine | Ave | 2.018 | 1.970 | | 9.76 | 10.0 | -2.4 | 30.0 |
| Benzaldehyde | Ave | 1.420 | 1.308 | 0.0100 | 18.4 | 20.0 | -7.9 | 30.0 |
| Phenol | Ave | 2.099 | 2.079 | 0.8000 | 9.91 | 10.0 | -0.9 | 30.0 |
| Aniline | Ave | 2.617 | 2.563 | | 9.79 | 10.0 | -2.1 | 30.0 |
| Bis(2-chloroethyl)ether | Ave | 1.803 | 1.702 | 0.7000 | 9.44 | 10.0 | -5.6 | 30.0 |
| 2-Chlorophenol | Ave | 1.517 | 1.526 | 0.8000 | 10.1 | 10.0 | 0.6 | 30.0 |
| n-Decane | Ave | 1.758 | 1.658 | | 9.44 | 10.0 | -5.6 | 30.0 |
| 1,3-Dichlorobenzene | Ave | 1.573 | 1.505 | | 9.57 | 10.0 | -4.3 | 30.0 |
| 1,4-Dichlorobenzene | Ave | 1.602 | 1.552 | | 9.69 | 10.0 | -3.1 | 30.0 |
| Benzyl alcohol | Ave | 1.029 | 1.068 | | 10.4 | 10.0 | 3.8 | 30.0 |
| 1,2-Dichlorobenzene | Ave | 1.540 | 1.462 | | 9.49 | 10.0 | -5.1 | 30.0 |
| 2-Methylphenol | Ave | 1.570 | 1.558 | 0.7000 | 9.93 | 10.0 | -0.7 | 30.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.024 | 1.968 | | 9.73 | 10.0 | -2.7 | 30.0 |
| Indene | Ave | 2.688 | 2.473 | | 18.4 | 20.0 | -8.0 | 30.0 |
| 3 & 4 Methylphenol | Ave | 1.620 | 1.612 | | 9.96 | 10.0 | -0.4 | 30.0 |
| N-Nitrosodi-n-propylamine | Ave | 1.139 | 1.142 | 0.5000 | 10.0 | 10.0 | 0.3 | 30.0 |
| Acetophenone | Ave | 2.353 | 2.256 | 0.0100 | 9.59 | 10.0 | -4.1 | 30.0 |
| Hexachloroethane | Ave | 0.6683 | 0.6557 | 0.3000 | 9.81 | 10.0 | -1.9 | 30.0 |
| Nitrobenzene | Ave | 0.4459 | 0.4395 | 0.2000 | 9.86 | 10.0 | -1.4 | 30.0 |
| Isophorone | Ave | 0.7626 | 0.7927 | 0.4000 | 10.4 | 10.0 | 3.9 | 30.0 |
| 2-Nitrophenol | Ave | 0.1660 | 0.1808 | 0.1000 | 10.9 | 10.0 | 8.9 | 30.0 |
| 2,4-Dimethylphenol | Ave | 0.3818 | 0.3725 | 0.2000 | 9.76 | 10.0 | -2.4 | 30.0 |
| Benzoic acid | Lin1 | | 0.2152 | | 20.9 | 20.0 | 4.4 | 30.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4539 | 0.4473 | 0.3000 | 9.86 | 10.0 | -1.4 | 30.0 |
| 2,4-Dichlorophenol | Ave | 0.2489 | 0.2521 | 0.2000 | 10.1 | 10.0 | 1.3 | 30.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.2911 | 0.2788 | | 9.58 | 10.0 | -4.2 | 30.0 |
| Naphthalene | Ave | 1.093 | 1.042 | 0.7000 | 9.53 | 10.0 | -4.7 | 30.0 |
| 4-Chloroaniline | Ave | 0.4651 | 0.4784 | 0.0100 | 10.3 | 10.0 | 2.9 | 30.0 |
| 2,6-Dichlorophenol | Ave | 0.2535 | 0.2551 | | 10.1 | 10.0 | 0.6 | 30.0 |
| Hexachlorobutadiene | Ave | 0.1548 | 0.1494 | 0.0100 | 9.65 | 10.0 | -3.5 | 30.0 |
| Caprolactam | Ave | 0.1118 | 0.1204 | 0.0100 | 21.5 | 20.0 | 7.7 | 30.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3177 | 0.3422 | 0.2000 | 10.8 | 10.0 | 7.7 | 30.0 |
| 2-Methylnaphthalene | Ave | 0.6768 | 0.6510 | 0.4000 | 9.62 | 10.0 | -3.8 | 30.0 |
| 1-Methylnaphthalene | Ave | 0.6004 | 0.6115 | | 10.2 | 10.0 | 1.9 | 30.0 |
| Hexachlorocyclopentadiene | Ave | 0.2585 | 0.2932 | 0.0500 | 11.3 | 10.0 | 13.4 | 30.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4935 | 0.4572 | 0.0100 | 9.26 | 10.0 | -7.4 | 30.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3377 | 0.3579 | 0.2000 | 10.6 | 10.0 | 6.0 | 30.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3511 | 0.3667 | 0.2000 | 10.4 | 10.0 | 4.4 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204569/11 Calibration Date: 10/30/2015 17:08

Instrument ID: A4HP9 Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS ID: 0.45(mm) Calib End Date: 10/30/2015 16:42

Lab File ID: 51030011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 1,1'-Biphenyl | Ave | 1.542 | 1.453 | 0.0100 | 9.42 | 10.0 | -5.8 | 30.0 |
| 2-Chloronaphthalene | Ave | 1.162 | 1.117 | 0.8000 | 9.62 | 10.0 | -3.8 | 30.0 |
| 2-Nitroaniline | Ave | 0.4031 | 0.4426 | 0.0100 | 11.0 | 10.0 | 9.8 | 30.0 |
| Dimethyl phthalate | Ave | 1.213 | 1.228 | 0.0100 | 10.1 | 10.0 | 1.2 | 30.0 |
| 1,3-Dinitrobenzene | Ave | 0.1896 | 0.2188 | | 11.5 | 10.0 | 15.4 | 30.0 |
| 2,6-Dinitrotoluene | Ave | 0.2814 | 0.2991 | | 10.6 | 10.0 | 6.3 | 30.0 |
| Acenaphthylene | Ave | 1.895 | 1.849 | 0.9000 | 9.76 | 10.0 | -2.4 | 30.0 |
| 3-Nitroaniline | Ave | 0.3457 | 0.3664 | 0.0100 | 10.6 | 10.0 | 6.0 | 30.0 |
| 2,4-Dinitrophenol | Lin | | 0.1277 | 0.0100 | 19.3 | 20.0 | -3.4 | 30.0 |
| Acenaphthene | Ave | 1.225 | 1.185 | 0.9000 | 9.67 | 10.0 | -3.3 | 30.0 |
| 4-Nitrophenol | Ave | 0.2054 | 0.2330 | | 22.7 | 20.0 | 13.4 | 30.0 |
| 2,4-Dinitrotoluene | Ave | 0.3558 | 0.3806 | 0.2000 | 10.7 | 10.0 | 7.0 | 30.0 |
| Dibenzofuran | Ave | 1.646 | 1.554 | 0.8000 | 9.44 | 10.0 | -5.6 | 30.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.2627 | 0.2874 | 0.0100 | 10.9 | 10.0 | 9.4 | 30.0 |
| Diethyl phthalate | Ave | 1.206 | 1.131 | 0.0100 | 9.38 | 10.0 | -6.2 | 30.0 |
| Hexadecane | Ave | 0.8758 | 0.8247 | | 9.42 | 10.0 | -5.8 | 30.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.5909 | 0.5890 | 0.4000 | 9.97 | 10.0 | -0.3 | 30.0 |
| 4-Nitroaniline | Ave | 0.3467 | 0.3816 | 0.0100 | 11.0 | 10.0 | 10.1 | 30.0 |
| Fluorene | Ave | 1.288 | 1.273 | 0.9000 | 9.89 | 10.0 | -1.1 | 30.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1284 | 0.0100 | 21.0 | 20.0 | 4.8 | 30.0 |
| N-Nitrosodiphenylamine | Ave | 0.5262 | 0.5266 | 0.0100 | 20.0 | 20.0 | 0.0 | 30.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 1.014 | 1.113 | | 11.0 | 10.0 | 9.8 | 30.0 |
| Azobenzene | Ave | 1.014 | 1.112 | | 11.0 | 10.0 | 9.7 | 30.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.1862 | 0.1987 | 0.1000 | 10.7 | 10.0 | 6.7 | 30.0 |
| Hexachlorobenzene | Ave | 0.1906 | 0.1906 | 0.1000 | 10.0 | 10.0 | 0.0 | 30.0 |
| Atrazine | Ave | 0.1757 | 0.1916 | 0.0100 | 21.8 | 20.0 | 9.0 | 30.0 |
| n-Octadecane | Ave | 0.5128 | 0.5469 | | 10.7 | 10.0 | 6.6 | 30.0 |
| Pentachlorophenol | Lin1 | | 0.1354 | 0.0500 | 22.2 | 20.0 | 11.1 | 30.0 |
| Phenanthrene | Ave | 1.179 | 1.105 | 0.7000 | 9.38 | 10.0 | -6.2 | 30.0 |
| Anthracene | Ave | 1.143 | 1.148 | 0.7000 | 10.0 | 10.0 | 0.4 | 30.0 |
| Carbazole | Ave | 0.996 | 1.069 | 0.0100 | 10.7 | 10.0 | 7.3 | 30.0 |
| Di-n-butyl phthalate | Ave | 1.181 | 1.389 | 0.0100 | 11.8 | 10.0 | 17.6 | 30.0 |
| Fluoranthene | Ave | 1.125 | 1.216 | 0.6000 | 10.8 | 10.0 | 8.0 | 30.0 |
| Benzidine | Lin1 | | 0.6541 | | 20.2 | 20.0 | 0.8 | 30.0 |
| Pyrene | Ave | 1.165 | 1.126 | 0.6000 | 9.66 | 10.0 | -3.4 | 30.0 |
| Butyl benzyl phthalate | Ave | 0.5273 | 0.5758 | 0.0100 | 10.9 | 10.0 | 9.2 | 30.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.7086 | 0.7773 | 0.0100 | 11.0 | 10.0 | 9.7 | 30.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3426 | 0.3569 | 0.0100 | 20.8 | 20.0 | 4.2 | 30.0 |
| Benzo[a]anthracene | Ave | 1.114 | 1.131 | 0.8000 | 10.2 | 10.0 | 1.6 | 30.0 |
| Chrysene | Ave | 1.127 | 1.120 | 0.7000 | 9.93 | 10.0 | -0.7 | 30.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-204569/11 Calibration Date: 10/30/2015 17:08

Instrument ID: A4HP9 Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS ID: 0.45(mm) Calib End Date: 10/30/2015 16:42

Lab File ID: 51030011.D Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Di-n-octyl phthalate | Lin1 | | 1.598 | 0.0100 | 10.3 | 10.0 | 2.7 | 30.0 |
| Benzo[b]fluoranthene | Ave | 1.034 | 1.155 | 0.7000 | 11.2 | 10.0 | 11.8 | 30.0 |
| Benzo[k]fluoranthene | Ave | 1.334 | 1.317 | 0.7000 | 9.87 | 10.0 | -1.3 | 30.0 |
| Benzo[a]pyrene | Ave | 1.098 | 1.155 | 0.7000 | 10.5 | 10.0 | 5.2 | 30.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.106 | 1.175 | 0.5000 | 10.6 | 10.0 | 6.2 | 30.0 |
| Dibenz(a,h)anthracene | Ave | 0.8758 | 0.9878 | 0.4000 | 11.3 | 10.0 | 12.8 | 30.0 |
| Benzo[g,h,i]perylene | Ave | 0.9937 | 1.012 | 0.5000 | 10.2 | 10.0 | 1.9 | 30.0 |
| 2-Fluorophenol (Surr) | Ave | 1.509 | 1.588 | | 10.5 | 10.0 | 5.2 | 30.0 |
| Phenol-d5 (Surr) | Ave | 1.961 | 2.063 | | 10.5 | 10.0 | 5.2 | 30.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.4689 | 0.4903 | | 10.5 | 10.0 | 4.6 | 30.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.261 | 1.273 | | 10.1 | 10.0 | 1.0 | 30.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1401 | 0.1609 | | 11.5 | 10.0 | 14.8 | 30.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7313 | 0.7754 | | 10.6 | 10.0 | 6.0 | 30.0 |

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab File ID: 51111001.D  DFTPP Injection Date: 11/11/2015

Instrument ID: A4HP9  DFTPP Injection Time: 07:27

Analysis Batch No.: 206233

| M/E | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE | |
|---|---|---|---|
| 51 | 30.0- 80.0% of mass 198 | 50.3 | |
| 68 | Less than 2.0% of mass 69 | 0.0 | (0.0)1 |
| 69 | Mass 69 relative abundance | 58.1 | |
| 70 | Less than 2.0% of mass 69 | 0.3 | (0.5)1 |
| 127 | 25.0 - 75.0% of mass 198 | 54.3 | |
| 197 | Less than 1.0% of mass 198 | 0.0 | |
| 198 | Base Peak, 100% relative abundance | 100.0 | |
| 199 | 5.0 to 9.0% of mass 198 | 6.6 | |
| 275 | 10.0- 30.0% of mass 198 | 22.9 | |
| 365 | Greater than 0.75% of mass 198 | 2.7 | |
| 441 | Present, but less than mass 443 | 10.7 | |
| 442 | 40.0 - 110.0% of mass 198 | 61.4 | |
| 443 | 15.0 - 24.0% of mass 442 | 11.9 | (19.4)2 |

1-Value is % mass 69  2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

| CLIENT SAMPLE ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|---|---|---|---|---|
| | CCV 240-206233/2 | 51111002.D | 11/11/2015 | 07:49 |
| | MB 240-205866/8-A | 51111006.D | 11/11/2015 | 09:54 |
| | LCS 240-205866/9-A | 51111007.D | 11/11/2015 | 10:19 |
| F-VB-UST-11-14 | 240-57597-11 | 51111019.D | 11/11/2015 | 14:50 |
| F-VB-UST-12-14 | 240-57597-12 | 51111020.D | 11/11/2015 | 15:14 |
| F-VB-UST-13-14 | 240-57597-13 | 51111021.D | 11/11/2015 | 15:39 |
| F-VB-UST-14-14 | 240-57597-14 | 51111022.D | 11/11/2015 | 16:03 |
| F-VB-UST-DUP | 240-57597-15 | 51111023.D | 11/11/2015 | 16:28 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 51111024.D | 11/11/2015 | 16:52 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 51111025.D | 11/11/2015 | 17:17 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206233/2  Calibration Date: 11/11/2015 07:49

Instrument ID: A4HP9  Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS  ID: 0.45(mm)  Calib End Date: 10/30/2015 16:42

Lab File ID: 51111002.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| N-Nitrosodimethylamine | Ave | 1.097 | 1.401 | | 12.8 | 10.0 | 27.7* | 20.0 |
| Pyridine | Ave | 2.018 | 2.368 | | 11.7 | 10.0 | 17.4 | 20.0 |
| Benzaldehyde | Ave | 1.420 | 0.9711 | 0.0100 | 13.7 | 20.0 | -31.6* | 20.0 |
| Phenol | Ave | 2.099 | 2.304 | 0.8000 | 11.0 | 10.0 | 9.8 | 20.0 |
| Aniline | Ave | 2.617 | 2.827 | | 10.8 | 10.0 | 8.0 | 20.0 |
| Bis(2-chloroethyl)ether | Ave | 1.803 | 2.052 | 0.7000 | 11.4 | 10.0 | 13.8 | 20.0 |
| 2-Chlorophenol | Ave | 1.517 | 1.503 | 0.8000 | 9.91 | 10.0 | -0.9 | 20.0 |
| n-Decane | Ave | 1.758 | 2.188 | | 12.4 | 10.0 | 24.5* | 20.0 |
| 1,3-Dichlorobenzene | Ave | 1.573 | 1.567 | | 9.97 | 10.0 | -0.3 | 20.0 |
| 1,4-Dichlorobenzene | Ave | 1.602 | 1.600 | | 9.99 | 10.0 | -0.1 | 20.0 |
| Benzyl alcohol | Ave | 1.029 | 1.117 | | 10.9 | 10.0 | 8.6 | 20.0 |
| 1,2-Dichlorobenzene | Ave | 1.540 | 1.527 | | 9.91 | 10.0 | -0.9 | 20.0 |
| 2-Methylphenol | Ave | 1.570 | 1.556 | 0.7000 | 9.91 | 10.0 | -0.9 | 20.0 |
| bis (2-chloroisopropyl) ether | Ave | 2.024 | 2.359 | | 11.7 | 10.0 | 16.6 | 20.0 |
| Indene | Ave | 2.688 | 2.657 | | 19.8 | 20.0 | -1.1 | 20.0 |
| 3 & 4 Methylphenol | Ave | 1.620 | 1.541 | | 9.51 | 10.0 | -4.9 | 20.0 |
| N-Nitrosodi-n-propylamine | Ave | 1.139 | 1.450 | 0.5000 | 12.7 | 10.0 | 27.4* | 20.0 |
| Acetophenone | Ave | 2.353 | 2.244 | 0.0100 | 9.54 | 10.0 | -4.6 | 20.0 |
| Hexachloroethane | Ave | 0.6683 | 0.6695 | 0.3000 | 10.0 | 10.0 | 0.2 | 20.0 |
| Nitrobenzene | Ave | 0.4459 | 0.4894 | 0.2000 | 11.0 | 10.0 | 9.8 | 20.0 |
| Isophorone | Ave | 0.7626 | 0.8151 | 0.4000 | 10.7 | 10.0 | 6.9 | 20.0 |
| 2-Nitrophenol | Ave | 0.1660 | 0.1725 | 0.1000 | 10.4 | 10.0 | 3.9 | 20.0 |
| 2,4-Dimethylphenol | Ave | 0.3818 | 0.3609 | 0.2000 | 9.45 | 10.0 | -5.5 | 20.0 |
| Benzoic acid | Lin1 | | 0.1419 | | 14.3 | 20.0 | -28.4* | 20.0 |
| Bis(2-chloroethoxy)methane | Ave | 0.4539 | 0.4711 | 0.3000 | 10.4 | 10.0 | 3.8 | 20.0 |
| 2,4-Dichlorophenol | Ave | 0.2489 | 0.2500 | 0.2000 | 10.0 | 10.0 | 0.4 | 20.0 |
| 1,2,4-Trichlorobenzene | Ave | 0.2911 | 0.2488 | | 8.54 | 10.0 | -14.6 | 20.0 |
| Naphthalene | Ave | 1.093 | 0.9231 | 0.7000 | 8.44 | 10.0 | -15.6 | 20.0 |
| 4-Chloroaniline | Ave | 0.4651 | 0.3947 | 0.0100 | 8.49 | 10.0 | -15.1 | 20.0 |
| 2,6-Dichlorophenol | Ave | 0.2535 | 0.2456 | | 9.69 | 10.0 | -3.1 | 20.0 |
| Hexachlorobutadiene | Ave | 0.1548 | 0.1309 | 0.0100 | 8.46 | 10.0 | -15.4 | 20.0 |
| Caprolactam | Ave | 0.1118 | 0.1051 | 0.0100 | 18.8 | 20.0 | -6.0 | 20.0 |
| 4-Chloro-3-methylphenol | Ave | 0.3177 | 0.3036 | 0.2000 | 9.56 | 10.0 | -4.4 | 20.0 |
| 2-Methylnaphthalene | Ave | 0.6768 | 0.6149 | 0.4000 | 9.09 | 10.0 | -9.1 | 20.0 |
| 1-Methylnaphthalene | Ave | 0.6004 | 0.5452 | | 9.08 | 10.0 | -9.2 | 20.0 |
| Hexachlorocyclopentadiene | Ave | 0.2585 | 0.2068 | 0.0500 | 8.00 | 10.0 | -20.0 | 20.0 |
| 1,2,4,5-Tetrachlorobenzene | Ave | 0.4935 | 0.4285 | 0.0100 | 8.68 | 10.0 | -13.2 | 20.0 |
| 2,4,6-Trichlorophenol | Ave | 0.3377 | 0.3103 | 0.2000 | 9.19 | 10.0 | -8.1 | 20.0 |
| 2,4,5-Trichlorophenol | Ave | 0.3511 | 0.3097 | 0.2000 | 8.82 | 10.0 | -11.8 | 20.0 |
| 1,1'-Biphenyl | Ave | 1.542 | 1.307 | 0.0100 | 8.48 | 10.0 | -15.2 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206233/2  Calibration Date: 11/11/2015 07:49

Instrument ID: A4HP9  Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS  ID: 0.45(mm)  Calib End Date: 10/30/2015 16:42

Lab File ID: 51111002.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| 2-Chloronaphthalene | Ave | 1.162 | 1.068 | 0.8000 | 9.19 | 10.0 | -8.1 | 20.0 |
| 2-Nitroaniline | Ave | 0.4031 | 0.4647 | 0.0100 | 11.5 | 10.0 | 15.3 | 20.0 |
| Dimethyl phthalate | Ave | 1.213 | 1.112 | 0.0100 | 9.17 | 10.0 | -8.3 | 20.0 |
| 1,3-Dinitrobenzene | Ave | 0.1896 | 0.2036 | | 10.7 | 10.0 | 7.4 | 20.0 |
| 2,6-Dinitrotoluene | Ave | 0.2814 | 0.2741 | | 9.74 | 10.0 | -2.6 | 20.0 |
| Acenaphthylene | Ave | 1.895 | 1.621 | 0.9000 | 8.55 | 10.0 | -14.5 | 20.0 |
| 3-Nitroaniline | Ave | 0.3457 | 0.3215 | 0.0100 | 9.30 | 10.0 | -7.0 | 20.0 |
| 2,4-Dinitrophenol | Lin | | 0.1118 | 0.0100 | 17.4 | 20.0 | -13.2 | 20.0 |
| Acenaphthene | Ave | 1.225 | 1.006 | 0.9000 | 8.21 | 10.0 | -17.9 | 20.0 |
| 4-Nitrophenol | Ave | 0.2054 | 0.1907 | | 18.6 | 20.0 | -7.2 | 20.0 |
| 2,4-Dinitrotoluene | Ave | 0.3558 | 0.3600 | 0.2000 | 10.1 | 10.0 | 1.2 | 20.0 |
| Dibenzofuran | Ave | 1.646 | 1.411 | 0.8000 | 8.57 | 10.0 | -14.3 | 20.0 |
| 2,3,4,6-Tetrachlorophenol | Ave | 0.2627 | 0.2179 | 0.0100 | 8.29 | 10.0 | -17.1 | 20.0 |
| Diethyl phthalate | Ave | 1.206 | 0.9799 | 0.0100 | 8.12 | 10.0 | -18.8 | 20.0 |
| Hexadecane | Ave | 0.8758 | 0.8735 | | 9.97 | 10.0 | -0.3 | 20.0 |
| 4-Chlorophenyl phenyl ether | Ave | 0.5909 | 0.5060 | 0.4000 | 8.56 | 10.0 | -14.4 | 20.0 |
| 4-Nitroaniline | Ave | 0.3467 | 0.3204 | 0.0100 | 9.24 | 10.0 | -7.6 | 20.0 |
| Fluorene | Ave | 1.288 | 1.141 | 0.9000 | 8.86 | 10.0 | -11.4 | 20.0 |
| 4,6-Dinitro-2-methylphenol | Lin1 | | 0.1238 | 0.0100 | 20.3 | 20.0 | 1.4 | 20.0 |
| N-Nitrosodiphenylamine | Ave | 0.5262 | 0.5308 | 0.0100 | 20.2 | 20.0 | 0.9 | 20.0 |
| 1,2-Diphenylhydrazine(as Azobenzene) | Ave | 1.014 | 1.110 | | 10.9 | 10.0 | 9.5 | 20.0 |
| Azobenzene | Ave | 1.014 | 1.111 | | 11.0 | 10.0 | 9.5 | 20.0 |
| 4-Bromophenyl phenyl ether | Ave | 0.1862 | 0.1790 | 0.1000 | 9.61 | 10.0 | -3.9 | 20.0 |
| Hexachlorobenzene | Ave | 0.1906 | 0.1928 | 0.1000 | 10.1 | 10.0 | 1.2 | 20.0 |
| Atrazine | Ave | 0.1757 | 0.1904 | 0.0100 | 21.7 | 20.0 | 8.3 | 20.0 |
| n-Octadecane | Ave | 0.5128 | 0.6714 | | 13.1 | 10.0 | 30.9* | 20.0 |
| Pentachlorophenol | Lin1 | | 0.1117 | 0.0500 | 18.5 | 20.0 | -7.7 | 20.0 |
| Phenanthrene | Ave | 1.179 | 1.084 | 0.7000 | 9.19 | 10.0 | -8.1 | 20.0 |
| Anthracene | Ave | 1.143 | 1.091 | 0.7000 | 9.54 | 10.0 | -4.6 | 20.0 |
| Carbazole | Ave | 0.996 | 1.073 | 0.0100 | 10.8 | 10.0 | 7.7 | 20.0 |
| Di-n-butyl phthalate | Ave | 1.181 | 1.258 | 0.0100 | 10.6 | 10.0 | 6.5 | 20.0 |
| Fluoranthene | Ave | 1.125 | 1.144 | 0.6000 | 10.2 | 10.0 | 1.7 | 20.0 |
| Benzidine | Lin1 | | 0.4427 | | 13.9 | 20.0 | -30.7* | 20.0 |
| Pyrene | Ave | 1.165 | 1.127 | 0.6000 | 9.67 | 10.0 | -3.3 | 20.0 |
| Butyl benzyl phthalate | Ave | 0.5273 | 0.5343 | 0.0100 | 10.1 | 10.0 | 1.3 | 20.0 |
| Bis(2-ethylhexyl) phthalate | Ave | 0.7086 | 0.7187 | 0.0100 | 10.1 | 10.0 | 1.4 | 20.0 |
| 3,3'-Dichlorobenzidine | Ave | 0.3426 | 0.3520 | 0.0100 | 20.5 | 20.0 | 2.7 | 20.0 |
| Benzo[a]anthracene | Ave | 1.114 | 1.008 | 0.8000 | 9.05 | 10.0 | -9.5 | 20.0 |
| Chrysene | Ave | 1.127 | 0.9562 | 0.7000 | 8.48 | 10.0 | -15.2 | 20.0 |
| Di-n-octyl phthalate | Lin1 | | 1.464 | 0.0100 | 9.43 | 10.0 | -5.7 | 20.0 |

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206233/2  Calibration Date: 11/11/2015 07:49

Instrument ID: A4HP9  Calib Start Date: 10/30/2015 13:19

GC Column: RXI-5SILMS  ID: 0.45(mm)  Calib End Date: 10/30/2015 16:42

Lab File ID: 51111002.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Benzo[b]fluoranthene | Ave | 1.034 | 1.122 | 0.7000 | 10.9 | 10.0 | 8.6 | 20.0 |
| Benzo[k]fluoranthene | Ave | 1.334 | 1.254 | 0.7000 | 9.40 | 10.0 | -6.0 | 20.0 |
| Benzo[a]pyrene | Ave | 1.098 | 1.113 | 0.7000 | 10.1 | 10.0 | 1.3 | 20.0 |
| Indeno[1,2,3-cd]pyrene | Ave | 1.106 | 1.182 | 0.5000 | 10.7 | 10.0 | 6.8 | 20.0 |
| Dibenz(a,h)anthracene | Ave | 0.8758 | 0.9640 | 0.4000 | 11.0 | 10.0 | 10.1 | 20.0 |
| Benzo[g,h,i]perylene | Ave | 0.9937 | 0.9829 | 0.5000 | 9.89 | 10.0 | -1.1 | 20.0 |
| 1,4-Dioxane | Ave | 0.7606 | | | 0.000490 | 10.0 | | |
| 2-Fluorophenol (Surr) | Ave | 1.509 | 1.580 | | 10.5 | 10.0 | 4.7 | 20.0 |
| Phenol-d5 (Surr) | Ave | 1.961 | 2.103 | | 10.7 | 10.0 | 7.2 | 20.0 |
| Nitrobenzene-d5 (Surr) | Ave | 0.4689 | 0.4936 | | 10.5 | 10.0 | 5.3 | 20.0 |
| 2-Fluorobiphenyl (Surr) | Ave | 1.261 | 1.114 | | 8.84 | 10.0 | -11.6 | 20.0 |
| 2,4,6-Tribromophenol (Surr) | Ave | 0.1401 | 0.1104 | | 7.88 | 10.0 | -21.2* | 20.0 |
| Terphenyl-d14 (Surr) | Ave | 0.7313 | 0.7199 | | 9.84 | 10.0 | -1.6 | 20.0 |

# GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205786  Batch Start Date: 11/09/15 10:35  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/10/15 02:35

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exBENZALDEHYD 00031 | exBNASPIKE 00053 | exBNASURR W 00028 | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-E-1 | F-VB-UST-1-14 | 3540C, 8270D | T | 29.99 g | 2 mL | | | 1 mL | |
| 240-57597-D-2 | F-VB-UST-2-14 | 3540C, 8270D | T | 30.16 g | 2 mL | | | 1 mL | |
| 240-57597-E-3 | F-VB-UST-3-14 | 3540C, 8270D | T | 30.03 g | 2 mL | | | 1 mL | |
| 240-57597-E-4 | F-VB-UST-4-14 | 3540C, 8270D | T | 29.84 g | 2 mL | | | 1 mL | |
| 240-57597-E-5 | F-VB-UST-5-14 | 3540C, 8270D | T | 30.28 g | 2 mL | | | 1 mL | |
| 240-57597-E-6 | F-VB-UST-6-14 | 3540C, 8270D | T | 30.15 g | 2 mL | | | 1 mL | |
| 240-57597-E-7 | F-VB-UST-7-14 | 3540C, 8270D | T | 29.53 g | 2 mL | | | 1 mL | |
| 240-57597-D-8 | F-VB-UST-8-14 | 3540C, 8270D | T | 30.20 g | 2 mL | | | 1 mL | |
| 240-57597-E-9 | F-VB-UST-9-14 | 3540C, 8270D | T | 29.67 g | 2 mL | | | 1 mL | |
| 240-57597-E-10 | F-VB-UST-10-14 | 3540C, 8270D | T | 29.78 g | 2 mL | | | 1 mL | |
| MB 240-205786/23 | | 3540C, 8270D | | 30 g | 2 mL | | | 1 mL | |
| LCS 240-205786/24 | | 3540C, 8270D | | 30 g | 2 mL | 1 mL | 1 mL | 1 mL | |

| Batch Notes | |
|---|---|
| Balance ID | 1339912 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2258961 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205866  Batch Start Date: 11/09/15 01:00  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/10/15 05:00

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exBENZALDEHYD 00031 | exBNASPIKE 00053 | exBNASURR W 00028 | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-E-11 | F-VB-UST-11-14 | 3540C, 8270D | T | 29.86 g | 2 mL | | | 1 mL | |
| 240-57597-E-12 | F-VB-UST-12-14 | 3540C, 8270D | T | 29.84 g | 2 mL | | | 1 mL | |
| 240-57597-E-13 | F-VB-UST-13-14 | 3540C, 8270D | T | 30.37 g | 2 mL | | | 1 mL | |
| 240-57597-D-14 | F-VB-UST-14-14 | 3540C, 8270D | T | 29.64 g | 2 mL | | | 1 mL | |
| 240-57597-D-15 | F-VB-UST-DUP | 3540C, 8270D | T | 30.09 g | 2 mL | | | 1 mL | |
| 240-57597-D-15 MS | F-VB-UST-DUP | 3540C, 8270D | T | 30.05 g | 2 mL | 1 mL | 1 mL | 1 mL | |
| 240-57597-D-15 MSD | F-VB-UST-DUP | 3540C, 8270D | T | 29.93 g | 2 mL | 1 mL | 1 mL | 1 mL | |
| MB 240-205866/8 | | 3540C, 8270D | | 30 g | 2 mL | | | 1 mL | |
| LCS 240-205866/9 | | 3540C, 8270D | | 30 g | 2 mL | 1 mL | 1 mL | 1 mL | |

| Batch Notes | |
|---|---|
| Balance ID | 1339907 |
| Na2SO4 Lot Number | 2204509 |
| Prep Solvent Lot # | 2258960 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

FORM II
PCBS SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): CLP-2 ID: 0.53(mm)

| Client Sample ID | Lab Sample ID | TCX1 # | DCB1 # |
|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 78 | 87 |
| F-VB-UST-2-14 | 240-57597-2 | 84 | 100 |
| F-VB-UST-3-14 | 240-57597-3 | 68 | 95 |
| F-VB-UST-4-14 | 240-57597-4 | 75 | 104 |
| F-VB-UST-5-14 | 240-57597-5 | 73 | 91 |
| F-VB-UST-6-14 | 240-57597-6 | 76 | 109 |
| F-VB-UST-7-14 | 240-57597-7 | 76 | 98 |
| F-VB-UST-8-14 | 240-57597-8 | 72 | 90 |
| F-VB-UST-9-14 | 240-57597-9 | 77 | 95 |
| F-VB-UST-10-14 | 240-57597-10 | 83 | 89 |
| F-VB-UST-11-14 | 240-57597-11 | 82 | 95 |
| F-VB-UST-12-14 | 240-57597-12 | 81 | 92 |
| F-VB-UST-13-14 | 240-57597-13 | 77 | 94 |
| F-VB-UST-14-14 | 240-57597-14 | 78 | 89 |
| F-VB-UST-DUP | 240-57597-15 | 74 | 89 |
| | MB 240-205809/21-A | 75 | 89 |
| | LCS 240-205809/22-A | 75 | 85 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 81 | 95 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 77 | 89 |

| | QC LIMITS |
|---|---|
| TCX = Tetrachloro-m-xylene | 13-134 |
| DCB = DCB Decachlorobiphenyl | 10-155 |

# Column to be used to flag recovery values

FORM II 8082A

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: STD05 240-205222/31  Date Analyzed: 11/04/2015 22:13

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)

Lab File ID (Standard): P4000031.D  Heated Purge: (Y/N) N

Calibration ID: 31286

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| INITIAL CALIBRATION MID-POINT | | 138265983 | 2.08 | | | | |
| UPPER LIMIT | | 276531966 | 2.58 | | | | |
| LOWER LIMIT | | 69132992 | 1.58 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| ICV 240-205222/39 | | 151134439 | 2.09 | | | | |
| CCV 240-206283/3 CCVIS | | 120153607 | 2.07 | | | | |

BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206283/3  Date Analyzed: 11/11/2015 11:01

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)

Lab File ID (Standard): P4000003.D  Heated Purge: (Y/N) N

Calibration ID: 31286

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 120153607 | 2.07 | | | | |
| UPPER LIMIT | | 240307214 | 2.57 | | | | |
| LOWER LIMIT | | 60076804 | 1.57 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| CCV 240-206283/4 | | 117662635 | 2.07 | | | | |
| CCV 240-206283/5 | | 128931083 | 2.08 | | | | |
| CCV 240-206283/6 | | 129660063 | 2.07 | | | | |
| CCV 240-206283/7 | | 124415610 | 2.07 | | | | |
| 240-57597-1 | F-VB-UST-1-14 | 140229279 | 2.08 | | | | |
| 240-57597-2 | F-VB-UST-2-14 | 131329155 | 2.07 | | | | |
| 240-57597-3 | F-VB-UST-3-14 | 142270888 | 2.07 | | | | |
| 240-57597-4 | F-VB-UST-4-14 | 139052581 | 2.07 | | | | |
| 240-57597-5 | F-VB-UST-5-14 | 150647200 | 2.07 | | | | |
| 240-57597-6 | F-VB-UST-6-14 | 136592279 | 2.07 | | | | |
| 240-57597-7 | F-VB-UST-7-14 | 142766511 | 2.07 | | | | |
| 240-57597-8 | F-VB-UST-8-14 | 143364145 | 2.07 | | | | |
| 240-57597-9 | F-VB-UST-9-14 | 135564000 | 2.07 | | | | |
| 240-57597-10 | F-VB-UST-10-14 | 134509895 | 2.07 | | | | |
| 240-57597-11 | F-VB-UST-11-14 | 146836344 | 2.07 | | | | |
| 240-57597-12 | F-VB-UST-12-14 | 141041341 | 2.07 | | | | |
| 240-57597-13 | F-VB-UST-13-14 | 143471346 | 2.08 | | | | |
| 240-57597-14 | F-VB-UST-14-14 | 143338015 | 2.07 | | | | |
| MB 240-205809/21-A | | 150167483 | 2.07 | | | | |
| LCS 240-205809/22-A | | 138157065 | 2.07 | | | | |

BNB = 1-Bromo-2-nitrobenzene
BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM VIII
PCBS INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVIS 240-206283/28  Date Analyzed: 11/11/2015 17:50

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)

Lab File ID (Standard): P4000028.D  Heated Purge: (Y/N) N

Calibration ID: 31286

| | | BNB | | | | | |
|---|---|---|---|---|---|---|---|
| | | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 12/24 HOUR STD | | 124491700 | 2.08 | | | | |
| UPPER LIMIT | | 248983400 | 2.58 | | | | |
| LOWER LIMIT | | 62245850 | 1.58 | | | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | | | | | | |
| 240-57597-15 | F-VB-UST-DUP | 155311076 | 2.08 | | | | |
| 240-57597-15 MS | F-VB-UST-DUP MS | 144577114 | 2.07 | | | | |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 162028723 | 2.07 | | | | |

BNB = 1-Bromo-2-nitrobenzene

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

FORM VIII 8082A

FORM IV
PCBS METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: MB 240-205809/21-A

Matrix: Solid  Date Extracted: 11/09/2015 09:53

Lab File ID:(1) P4000026.D  Lab File ID:(2)

Date Analyzed:(1) 11/11/2015 17:17  Date Analyzed:(2)

Instrument ID:(1) A2HP4  Instrument ID:(2)

GC Column:(1) CLP-2  ID: 0.53(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 11/11/2015 13:28 | |
| F-VB-UST-2-14 | 240-57597-2 | 11/11/2015 13:45 | |
| F-VB-UST-3-14 | 240-57597-3 | 11/11/2015 14:01 | |
| F-VB-UST-4-14 | 240-57597-4 | 11/11/2015 14:17 | |
| F-VB-UST-5-14 | 240-57597-5 | 11/11/2015 14:33 | |
| F-VB-UST-6-14 | 240-57597-6 | 11/11/2015 14:50 | |
| F-VB-UST-7-14 | 240-57597-7 | 11/11/2015 15:06 | |
| F-VB-UST-8-14 | 240-57597-8 | 11/11/2015 15:23 | |
| F-VB-UST-9-14 | 240-57597-9 | 11/11/2015 15:39 | |
| F-VB-UST-10-14 | 240-57597-10 | 11/11/2015 15:55 | |
| F-VB-UST-11-14 | 240-57597-11 | 11/11/2015 16:12 | |
| F-VB-UST-12-14 | 240-57597-12 | 11/11/2015 16:28 | |
| F-VB-UST-13-14 | 240-57597-13 | 11/11/2015 16:44 | |
| F-VB-UST-14-14 | 240-57597-14 | 11/11/2015 17:01 | |
| | LCS 240-205809/22-A | 11/11/2015 17:33 | |
| F-VB-UST-DUP | 240-57597-15 | 11/11/2015 18:06 | |
| F-VB-UST-DUP MS | 240-57597-15 MS | 11/11/2015 18:22 | |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 11/11/2015 18:39 | |

FORM I
PCBS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205809/21-A

Matrix: Solid | Lab File ID: P4000026.D

Analysis Method: 8082A | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:53

Sample wt/vol: 30(g) | Date Analyzed: 11/11/2015 17:17

Con. Extract Vol.: 10(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: CLP-2 ID: 0.53(mm)

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206283 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 12674-11-2 | Aroclor-1016 | 12 | U | 33 | 12 |
| 11104-28-2 | Aroclor-1221 | 16 | U | 33 | 16 |
| 11141-16-5 | Aroclor-1232 | 20 | U | 33 | 20 |
| 53469-21-9 | Aroclor-1242 | 11 | U | 33 | 11 |
| 12672-29-6 | Aroclor-1248 | 8.0 | U | 33 | 8.0 |
| 11097-69-1 | Aroclor-1254 | 14 | U | 33 | 14 |
| 11096-82-5 | Aroclor-1260 | 9.0 | U | 33 | 9.0 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 877-09-8 | Tetrachloro-m-xylene | 75 | | 13-134 |
| 2051-24-3 | DCB Decachlorobiphenyl | 89 | | 10-155 |

FORM III
PCBS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P4000027.D

Lab ID: LCS 240-205809/22-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Aroclor-1016 | 333 | 196 | 59 | 51-120 | |
| Aroclor-1260 | 334 | 201 | 60 | 48-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM III
PCBS MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P4000030.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Aroclor-1016 | 394 | 14 U | 244 | 62 | 17-141 | |
| Aroclor-1260 | 395 | 10 U | 260 | 66 | 19-137 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

FORM III
PCBS MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: P4000031.D

Lab ID: 240-57597-15 MSD  Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Aroclor-1016 | 393 | 229 | 58 | 9 | 56 | 17-141 | |
| Aroclor-1260 | 393 | 238 | 61 | 9 | 55 | 19-137 | |

# Column to be used to flag recovery and RPD values

FORM III 8082A

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4 Start Date: 11/04/2015 14:51

Analysis Batch Number: 205222 End Date: 11/05/2015 00:57

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD005 240-205222/4 IC | | 11/04/2015 14:51 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/4 IC | | 11/04/2015 14:51 | 1 | P4000004.D | CLP-2 0.53(mm) |
| STD01 240-205222/5 IC | | 11/04/2015 15:08 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/5 IC | | 11/04/2015 15:08 | 1 | P4000005.D | CLP-2 0.53(mm) |
| STD02 240-205222/6 IC | | 11/04/2015 15:24 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/6 IC | | 11/04/2015 15:24 | 1 | P4000006.D | CLP-2 0.53(mm) |
| STD05 240-205222/7 IC | | 11/04/2015 15:40 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/7 IC | | 11/04/2015 15:40 | 1 | P4000007.D | CLP-2 0.53(mm) |
| STD1 240-205222/8 IC | | 11/04/2015 15:57 | 1 | | CLP-1 0.53(mm) |
| STD1 240-205222/8 IC | | 11/04/2015 15:57 | 1 | P4000008.D | CLP-2 0.53(mm) |
| STD2 240-205222/9 IC | | 11/04/2015 16:13 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/9 IC | | 11/04/2015 16:13 | 1 | P4000009.D | CLP-2 0.53(mm) |
| STD005 240-205222/10 IC | | 11/04/2015 16:30 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/10 IC | | 11/04/2015 16:30 | 1 | P4000010.D | CLP-2 0.53(mm) |
| STD01 240-205222/11 IC | | 11/04/2015 16:46 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/11 IC | | 11/04/2015 16:46 | 1 | P4000011.D | CLP-2 0.53(mm) |
| STD02 240-205222/12 IC | | 11/04/2015 17:02 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/12 IC | | 11/04/2015 17:02 | 1 | P4000012.D | CLP-2 0.53(mm) |
| STD05 240-205222/13 IC | | 11/04/2015 17:19 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/13 IC | | 11/04/2015 17:19 | 1 | P4000013.D | CLP-2 0.53(mm) |
| STD1 240-205222/14 IC | | 11/04/2015 17:35 | 1 | | CLP-1 0.53(mm) |
| STD1 240-205222/14 IC | | 11/04/2015 17:35 | 1 | P4000014.D | CLP-2 0.53(mm) |
| STD2 240-205222/15 IC | | 11/04/2015 17:51 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/15 IC | | 11/04/2015 17:51 | 1 | P4000015.D | CLP-2 0.53(mm) |
| STD005 240-205222/16 IC | | 11/04/2015 18:08 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/16 IC | | 11/04/2015 18:08 | 1 | P4000016.D | CLP-2 0.53(mm) |
| STD01 240-205222/17 IC | | 11/04/2015 18:24 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/17 IC | | 11/04/2015 18:24 | 1 | P4000017.D | CLP-2 0.53(mm) |
| STD02 240-205222/18 IC | | 11/04/2015 18:41 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/18 IC | | 11/04/2015 18:41 | 1 | P4000018.D | CLP-2 0.53(mm) |
| STD05 240-205222/19 IC | | 11/04/2015 18:57 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/19 IC | | 11/04/2015 18:57 | 1 | P4000019.D | CLP-2 0.53(mm) |
| STD1 240-205222/20 IC | | 11/04/2015 19:13 | 1 | | CLP-1 0.53(mm) |

8082A

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4  Start Date: 11/04/2015 14:51

Analysis Batch Number: 205222  End Date: 11/05/2015 00:57

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-205222/20 IC | | 11/04/2015 19:13 | 1 | P4000020.D | CLP-2 0.53(mm) |
| STD2 240-205222/21 IC | | 11/04/2015 19:30 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/21 IC | | 11/04/2015 19:30 | 1 | P4000021.D | CLP-2 0.53(mm) |
| STD005 240-205222/22 IC | | 11/04/2015 19:46 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/22 IC | | 11/04/2015 19:46 | 1 | P4000022.D | CLP-2 0.53(mm) |
| STD01 240-205222/23 IC | | 11/04/2015 20:02 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/23 IC | | 11/04/2015 20:02 | 1 | P4000023.D | CLP-2 0.53(mm) |
| STD02 240-205222/24 IC | | 11/04/2015 20:19 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/24 IC | | 11/04/2015 20:19 | 1 | P4000024.D | CLP-2 0.53(mm) |
| STD05 240-205222/25 IC | | 11/04/2015 20:35 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/25 IC | | 11/04/2015 20:35 | 1 | P4000025.D | CLP-2 0.53(mm) |
| STD1 240-205222/26 IC | | 11/04/2015 20:51 | 1 | | CLP-1 0.53(mm) |
| STD1 240-205222/26 IC | | 11/04/2015 20:51 | 1 | P4000026.D | CLP-2 0.53(mm) |
| STD2 240-205222/27 IC | | 11/04/2015 21:08 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/27 IC | | 11/04/2015 21:08 | 1 | P4000027.D | CLP-2 0.53(mm) |
| STD005 240-205222/28 IC | | 11/04/2015 21:24 | 1 | | CLP-1 0.53(mm) |
| STD005 240-205222/28 IC | | 11/04/2015 21:24 | 1 | P4000028.D | CLP-2 0.53(mm) |
| STD01 240-205222/29 IC | | 11/04/2015 21:40 | 1 | | CLP-1 0.53(mm) |
| STD01 240-205222/29 IC | | 11/04/2015 21:40 | 1 | P4000029.D | CLP-2 0.53(mm) |
| STD02 240-205222/30 IC | | 11/04/2015 21:57 | 1 | | CLP-1 0.53(mm) |
| STD02 240-205222/30 IC | | 11/04/2015 21:57 | 1 | P4000030.D | CLP-2 0.53(mm) |
| STD05 240-205222/31 ICIS | | 11/04/2015 22:13 | 1 | | CLP-1 0.53(mm) |
| STD05 240-205222/31 ICIS | | 11/04/2015 22:13 | 1 | P4000031.D | CLP-2 0.53(mm) |
| STD1 240-205222/32 IC | | 11/04/2015 22:29 | 1 | | CLP-1 0.53(mm) |
| STD1 240-205222/32 IC | | 11/04/2015 22:29 | 1 | P4000032.D | CLP-2 0.53(mm) |
| STD2 240-205222/33 IC | | 11/04/2015 22:46 | 1 | | CLP-1 0.53(mm) |
| STD2 240-205222/33 IC | | 11/04/2015 22:46 | 1 | P4000033.D | CLP-2 0.53(mm) |
| ICV 240-205222/34 | | 11/04/2015 23:02 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/34 | | 11/04/2015 23:02 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/35 | | 11/04/2015 23:19 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/35 | | 11/04/2015 23:19 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/36 | | 11/04/2015 23:35 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/36 | | 11/04/2015 23:35 | 1 | | CLP-2 0.53(mm) |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4  Start Date: 11/04/2015 14:51

Analysis Batch Number: 205222  End Date: 11/05/2015 00:57

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| ICV 240-205222/37 | | 11/04/2015 23:51 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/37 | | 11/04/2015 23:51 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/38 | | 11/05/2015 00:08 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/38 | | 11/05/2015 00:08 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/39 | | 11/05/2015 00:24 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/39 | | 11/05/2015 00:24 | 1 | P4000039.D | CLP-2 0.53(mm) |
| ICV 240-205222/40 | | 11/05/2015 00:40 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/40 | | 11/05/2015 00:40 | 1 | | CLP-2 0.53(mm) |
| ICV 240-205222/41 | | 11/05/2015 00:57 | 1 | | CLP-1 0.53(mm) |
| ICV 240-205222/41 | | 11/05/2015 00:57 | 1 | | CLP-2 0.53(mm) |

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 14:51  Calibration End Date: 11/04/2015 16:13  Calibration ID: 31262

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/4 | P4000004.D |
| Level 2 | STD01 240-205222/5 | P4000005.D |
| Level 3 | STD02 240-205222/6 | P4000006.D |
| Level 4 | STD05 240-205222/7 | P4000007.D |
| Level 5 | STD1 240-205222/8 | P4000008.D |
| Level 6 | STD2 240-205222/9 | P4000009.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1232 Peak 1 | 0.0146<br>0.0122 | 0.0144 | 0.0134 | 0.0134 | 0.0122 | Lin1 | 0.0002 | 0.0122 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1232 Peak 2 | 0.0119<br>0.0097 | 0.0118 | 0.0110 | 0.0109 | 0.0098 | Lin1 | 0.0002 | 0.0098 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 3 | 0.0217<br>0.0205 | 0.0220 | 0.0209 | 0.0216 | 0.0202 | Lin1 | 0.0001 | 0.0205 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1232 Peak 4 | 0.0095<br>0.0081 | 0.0100 | 0.0094 | 0.0089 | 0.0081 | Lin1 | 0.0001 | 0.0081 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1232 Peak 5 | 0.0072<br>0.0065 | 0.0075 | 0.0069 | 0.0070 | 0.0066 | Lin1 | 0.0001 | 0.0066 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1262 Peak 1 | 0.0260<br>0.0242 | 0.0274 | 0.0250 | 0.0248 | 0.0248 | Lin1 | 0.0002 | 0.0244 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 2 | 0.0377<br>0.0377 | 0.0413 | 0.0382 | 0.0381 | 0.0383 | Lin1 | 0.0001 | 0.0379 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 3 | 0.0311<br>0.0318 | 0.0344 | 0.0320 | 0.0318 | 0.0323 | Lin1 | 0 | 0.0319 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 4 | 0.0698<br>0.0769 | 0.0788 | 0.0741 | 0.0751 | 0.0769 | Lin1 | 0 | 0.0769 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1262 Peak 5 | 0.0515<br>0.0538 | 0.0573 | 0.0535 | 0.0535 | 0.0541 | Lin1 | 0 | 0.0539 | | | | | | | 1.0000 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 16:30  Calibration End Date: 11/04/2015 17:51  Calibration ID: 31268

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/10 | P4000010.D |
| Level 2 | STD01 240-205222/11 | P4000011.D |
| Level 3 | STD02 240-205222/12 | P4000012.D |
| Level 4 | STD05 240-205222/13 | P4000013.D |
| Level 5 | STD1 240-205222/14 | P4000014.D |
| Level 6 | STD2 240-205222/15 | P4000015.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1242 Peak 1 | 0.0108 0.0082 | 0.0098 | 0.0093 | 0.0090 | 0.0085 | Lin1 | 0.0002 | 0.0083 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 2 | 0.0208 0.0159 | 0.0186 | 0.0181 | 0.0174 | 0.0164 | Lin1 | 0.0003 | 0.0160 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1242 Peak 3 | 0.0407 0.0367 | 0.0373 | 0.0369 | 0.0376 | 0.0368 | Lin1 | 0.0002 | 0.0367 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1242 Peak 4 | 0.0183 0.0142 | 0.0165 | 0.0153 | 0.0150 | 0.0144 | Lin1 | 0.0002 | 0.0142 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1242 Peak 5 | 0.0169 0.0109 | 0.0131 | 0.0124 | 0.0117 | 0.0111 | Lin1 | 0.0003 | 0.0108 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 1 | 0.1004 0.0910 | 0.0938 | 0.0918 | 0.0903 | 0.0899 | Lin1 | 0.0004 | 0.0902 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 2 | 0.0969 0.0876 | 0.0916 | 0.0907 | 0.0881 | 0.0870 | Lin1 | 0.0005 | 0.0871 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 3 | 0.0848 0.0764 | 0.0793 | 0.0781 | 0.0771 | 0.0756 | Lin1 | 0.0004 | 0.0759 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 4 | 0.0354 0.0308 | 0.0332 | 0.0323 | 0.0312 | 0.0308 | Lin1 | 0.0003 | 0.0306 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1268 Peak 5 | 0.2560 0.2513 | 0.2411 | 0.2398 | 0.2407 | 0.2403 | Lin1 | 0 | 0.2467 | | | | | | | 0.9990 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 18:08  Calibration End Date: 11/04/2015 19:30  Calibration ID: 31274

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/16 | P4000016.D |
| Level 2 | STD01 240-205222/17 | P4000017.D |
| Level 3 | STD02 240-205222/18 | P4000018.D |
| Level 4 | STD05 240-205222/19 | P4000019.D |
| Level 5 | STD1 240-205222/20 | P4000020.D |
| Level 6 | STD2 240-205222/21 | P4000021.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1248 Peak 1 | 0.0262 0.0244 | 0.0257 | 0.0252 | 0.0241 | 0.0238 | Lin1 | 0.0001 | 0.0241 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 2 | 0.0147 0.0119 | 0.0141 | 0.0133 | 0.0123 | 0.0118 | Lin1 | 0.0002 | 0.0118 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 3 | 0.0221 0.0182 | 0.0209 | 0.0200 | 0.0185 | 0.0178 | Lin1 | 0.0003 | 0.0180 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1248 Peak 4 | 0.0230 0.0230 | 0.0209 | 0.0242 | 0.0230 | 0.0224 | Lin1 | 0 | 0.0229 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1248 Peak 5 | 0.0111 0.0116 | 0.0121 | 0.0125 | 0.0118 | 0.0111 | Lin1 | 0 | 0.0115 | | | | | | | 0.9990 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4  GC Column: CLP-2  ID: 0.53(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 19:46  Calibration End Date: 11/04/2015 21:08  Calibration ID: 31280

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/22 | P4000022.D |
| Level 2 | STD01 240-205222/23 | P4000023.D |
| Level 3 | STD02 240-205222/24 | P4000024.D |
| Level 4 | STD05 240-205222/25 | P4000025.D |
| Level 5 | STD1 240-205222/26 | P4000026.D |
| Level 6 | STD2 240-205222/27 | P4000027.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1221 Peak 1 | 0.0082 0.0059 | 0.0071 | 0.0067 | 0.0065 | 0.0060 | Lin1 | 0.0001 | 0.0059 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1221 Peak 2 | 0.0056 0.0038 | 0.0048 | 0.0046 | 0.0044 | 0.0040 | Lin1 | 0.0001 | 0.0039 | | | | | | | 0.9980 | | 0.9900 |
| PCB-1221 Peak 3 | 0.0205 0.0147 | 0.0173 | 0.0165 | 0.0161 | 0.0151 | Lin1 | 0.0003 | 0.0148 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1254 Peak 1 | 0.0288 0.0211 | 0.0242 | 0.0224 | 0.0222 | 0.0212 | Lin1 | 0.0004 | 0.0209 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 2 | 0.0430 0.0354 | 0.0364 | 0.0351 | 0.0362 | 0.0354 | Lin1 | 0.0003 | 0.0352 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 3 | 0.0328 0.0255 | 0.0276 | 0.0256 | 0.0267 | 0.0261 | Lin1 | 0.0003 | 0.0255 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1254 Peak 4 | 0.0251 0.0215 | 0.0209 | 0.0202 | 0.0214 | 0.0214 | Lin1 | 0.0001 | 0.0213 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1254 Peak 5 | 0.0446 0.0346 | 0.0372 | 0.0350 | 0.0360 | 0.0351 | Lin1 | 0.0004 | 0.0345 | | | | | | | 1.0000 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
PCBS BY INTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205222

SDG No.:

Instrument ID: A2HP4 GC Column: CLP-2 ID: 0.53(mm) Heated Purge: (Y/N) N

Calibration Start Date: 11/04/2015 21:24 Calibration End Date: 11/04/2015 22:46 Calibration ID: 31286

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD005 240-205222/28 | P4000028.D |
| Level 2 | STD01 240-205222/29 | P4000029.D |
| Level 3 | STD02 240-205222/30 | P4000030.D |
| Level 4 | STD05 240-205222/31 | P4000031.D |
| Level 5 | STD1 240-205222/32 | P4000032.D |
| Level 6 | STD2 240-205222/33 | P4000033.D |

| ANALYTE | RRF | | | | | CURVE TYPE | COEFFICIENT | | | # | MIN RRF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 6 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | | B | M1 | M2 | | | | | | | | |
| PCB-1016 Peak 1 | 0.0178<br>0.0128 | 0.0153 | 0.0141 | 0.0136 | 0.0134 | Lin1 | 0.0003 | 0.0128 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 2 | 0.0346<br>0.0247 | 0.0296 | 0.0276 | 0.0265 | 0.0260 | Lin1 | 0.0005 | 0.0249 | | | | | | | 0.9990 | | 0.9900 |
| PCB-1016 Peak 3 | 0.0686<br>0.0581 | 0.0603 | 0.0585 | 0.0592 | 0.0594 | Lin1 | 0.0004 | 0.0582 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 4 | 0.0299<br>0.0224 | 0.0253 | 0.0239 | 0.0233 | 0.0230 | Lin1 | 0.0004 | 0.0224 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1016 Peak 5 | 0.0240<br>0.0168 | 0.0200 | 0.0185 | 0.0179 | 0.0174 | Lin1 | 0.0004 | 0.0168 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 1 | 0.0436<br>0.0313 | 0.0346 | 0.0370 | 0.0366 | 0.0324 | Lin1 | 0.0006 | 0.0318 | | | | | | | 0.9970 | | 0.9900 |
| PCB-1260 Peak 2 | 0.0509<br>0.0407 | 0.0443 | 0.0434 | 0.0430 | 0.0419 | Lin1 | 0.0005 | 0.0409 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 3 | 0.0730<br>0.0619 | 0.0645 | 0.0639 | 0.0642 | 0.0633 | Lin1 | 0.0005 | 0.0622 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 4 | 0.0854<br>0.0747 | 0.0754 | 0.0759 | 0.0763 | 0.0752 | Lin1 | 0.0004 | 0.0747 | | | | | | | 1.0000 | | 0.9900 |
| PCB-1260 Peak 5 | 0.0603<br>0.0505 | 0.0529 | 0.0527 | 0.0524 | 0.0514 | Lin1 | 0.0004 | 0.0506 | | | | | | | 1.0000 | | 0.9900 |
| Tetrachloro-m-xylene | 0.6148<br>0.6612 | 0.6230 | 0.6064 | 0.6560 | 0.6839 | Lin1 | 0 | 0.6687 | | | | | | | 0.9990 | | 0.9900 |
| DCB Decachlorobiphenyl | 0.6529<br>0.5587 | 0.6535 | 0.6238 | 0.6147 | 0.5968 | Lin1 | 0.0003 | 0.5721 | | | | | | | 0.9980 | | 0.9900 |

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-205222/39  Calibration Date: 11/05/2015 00:24

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 21:24

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 22:46

Lab File ID: P4000039.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0108 | | 0.822 | 1.00 | -17.8 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0217 | | 0.853 | 1.00 | -14.7 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0497 | | 0.848 | 1.00 | -15.2 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0191 | | 0.837 | 1.00 | -16.3 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0142 | | 0.824 | 1.00 | -17.6 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0258 | | 0.789 | 1.00 | -21.1* | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0324 | | 0.781 | 1.00 | -21.9* | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0430 | | 0.685 | 1.00 | -31.5* | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0708 | | 0.943 | 1.00 | -5.7 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0461 | | 0.903 | 1.00 | -9.7 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6638 | | 0.0500 | 0.0500 | -0.0 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.5091 | | 0.0439 | 0.0500 | -12.2 | 20.0 |

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4 Start Date: 11/11/2015 11:01

Analysis Batch Number: 206283 End Date: 11/11/2015 18:39

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-206283/3 CCVIS | | 11/11/2015 11:01 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/3 CCVIS | | 11/11/2015 11:01 | 1 | P4000003.D | CLP-2 0.53(mm) |
| CCV 240-206283/4 | | 11/11/2015 11:17 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/4 | | 11/11/2015 11:17 | 1 | P4000004.D | CLP-2 0.53(mm) |
| CCV 240-206283/5 | | 11/11/2015 11:34 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/5 | | 11/11/2015 11:34 | 1 | P4000005.D | CLP-2 0.53(mm) |
| CCV 240-206283/6 | | 11/11/2015 11:50 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/6 | | 11/11/2015 11:50 | 1 | P4000006.D | CLP-2 0.53(mm) |
| CCV 240-206283/7 | | 11/11/2015 12:06 | 1 | | CLP-1 0.53(mm) |
| CCV 240-206283/7 | | 11/11/2015 12:06 | 1 | P4000007.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 13:28 | 1 | | CLP-1 0.53(mm) |
| 240-57597-1 | F-VB-UST-1-14 | 11/11/2015 13:28 | 1 | P4000012.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 13:45 | 1 | | CLP-1 0.53(mm) |
| 240-57597-2 | F-VB-UST-2-14 | 11/11/2015 13:45 | 1 | P4000013.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 14:01 | 1 | | CLP-1 0.53(mm) |
| 240-57597-3 | F-VB-UST-3-14 | 11/11/2015 14:01 | 1 | P4000014.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 14:17 | 1 | | CLP-1 0.53(mm) |
| 240-57597-4 | F-VB-UST-4-14 | 11/11/2015 14:17 | 1 | P4000015.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 14:33 | 1 | | CLP-1 0.53(mm) |
| 240-57597-5 | F-VB-UST-5-14 | 11/11/2015 14:33 | 1 | P4000016.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 14:50 | 1 | | CLP-1 0.53(mm) |
| 240-57597-6 | F-VB-UST-6-14 | 11/11/2015 14:50 | 1 | P4000017.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 15:06 | 1 | | CLP-1 0.53(mm) |
| 240-57597-7 | F-VB-UST-7-14 | 11/11/2015 15:06 | 1 | P4000018.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 15:23 | 1 | | CLP-1 0.53(mm) |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 15:23 | 1 | P4000019.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 15:39 | 1 | | CLP-1 0.53(mm) |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 15:39 | 1 | P4000020.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 15:55 | 1 | | CLP-1 0.53(mm) |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 15:55 | 1 | P4000021.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 16:12 | 1 | | CLP-1 0.53(mm) |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 16:12 | 1 | P4000022.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 16:28 | 1 | | CLP-1 0.53(mm) |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 16:28 | 1 | P4000023.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 16:44 | 1 | | CLP-1 0.53(mm) |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 16:44 | 1 | P4000024.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 17:01 | 1 | | CLP-1 0.53(mm) |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 17:01 | 1 | P4000025.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 17:17 | 1 | | CLP-1 0.53(mm) |
| MB 240-205809/21-A | | 11/11/2015 17:17 | 1 | P4000026.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 17:33 | 1 | | CLP-1 0.53(mm) |
| LCS 240-205809/22-A | | 11/11/2015 17:33 | 1 | P4000027.D | CLP-2 0.53(mm) |
| CCVIS 240-206283/28 | | 11/11/2015 17:50 | 1 | | CLP-1 0.53(mm) |
| CCVIS 240-206283/28 | | 11/11/2015 17:50 | 1 | P4000028.D | CLP-2 0.53(mm) |

8082A

PCBS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP4  Start Date: 11/11/2015 11:01

Analysis Batch Number: 206283  End Date: 11/11/2015 18:39

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| ZZZZZ | | 11/11/2015 18:06 | 1 | | CLP-1 0.53(mm) |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 18:06 | 1 | P4000029.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 18:22 | 1 | | CLP-1 0.53(mm) |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 18:22 | 1 | P4000030.D | CLP-2 0.53(mm) |
| ZZZZZ | | 11/11/2015 18:39 | 1 | | CLP-1 0.53(mm) |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 18:39 | 1 | P4000031.D | CLP-2 0.53(mm) |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/3  Calibration Date: 11/11/2015 11:01

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 21:24

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 22:46

Lab File ID: P4000003.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0146 | | 0.549 | 0.500 | 9.8 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0282 | | 0.545 | 0.500 | 9.0 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0618 | | 0.525 | 0.500 | 5.0 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0246 | | 0.534 | 0.500 | 6.7 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0186 | | 0.530 | 0.500 | 6.1 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0333 | | 0.503 | 0.500 | 0.7 | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0427 | | 0.510 | 0.500 | 1.9 | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0632 | | 0.501 | 0.500 | 0.3 | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0741 | | 0.491 | 0.500 | -1.8 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0509 | | 0.494 | 0.500 | -1.2 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6986 | | 0.0264 | 0.0250 | 5.8 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.5950 | | 0.0254 | 0.0250 | 1.6 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/4  Calibration Date: 11/11/2015 11:17

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 14:51

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 16:13

Lab File ID: P4000004.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1232 Peak 1 | Lin1 | | 0.0149 | | 0.593 | 0.500 | 18.6 | 20.0 |
| PCB-1232 Peak 2 | Lin1 | | 0.0121 | | 0.604 | 0.500 | 20.8* | 20.0 |
| PCB-1232 Peak 3 | Lin1 | | 0.0235 | | 0.569 | 0.500 | 13.8 | 20.0 |
| PCB-1232 Peak 4 | Lin1 | | 0.0099 | | 0.592 | 0.500 | 18.4 | 20.0 |
| PCB-1232 Peak 5 | Lin1 | | 0.0080 | | 0.600 | 0.500 | 19.9 | 20.0 |
| PCB-1262 Peak 1 | Lin1 | | 0.0266 | | 0.539 | 0.500 | 7.8 | 20.0 |
| PCB-1262 Peak 2 | Lin1 | | 0.0407 | | 0.535 | 0.500 | 7.0 | 20.0 |
| PCB-1262 Peak 3 | Lin1 | | 0.0338 | | 0.527 | 0.500 | 5.5 | 20.0 |
| PCB-1262 Peak 4 | Lin1 | | 0.0787 | | 0.515 | 0.500 | 3.0 | 20.0 |
| PCB-1262 Peak 5 | Lin1 | | 0.0557 | | 0.517 | 0.500 | 3.3 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/5  Calibration Date: 11/11/2015 11:34

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 16:30

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 17:51

Lab File ID: P4000005.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1242 Peak 1 | Lin1 | | 0.0100 | | 0.583 | 0.500 | 16.7 | 20.0 |
| PCB-1242 Peak 2 | Lin1 | | 0.0189 | | 0.571 | 0.500 | 14.1 | 20.0 |
| PCB-1242 Peak 3 | Lin1 | | 0.0402 | | 0.543 | 0.500 | 8.7 | 20.0 |
| PCB-1242 Peak 4 | Lin1 | | 0.0163 | | 0.557 | 0.500 | 11.4 | 20.0 |
| PCB-1242 Peak 5 | Lin1 | | 0.0127 | | 0.562 | 0.500 | 12.3 | 20.0 |
| PCB-1268 Peak 1 | Lin1 | | 0.0893 | | 0.490 | 0.500 | -1.9 | 20.0 |
| PCB-1268 Peak 2 | Lin1 | | 0.0866 | | 0.491 | 0.500 | -1.8 | 20.0 |
| PCB-1268 Peak 3 | Lin1 | | 0.0761 | | 0.496 | 0.500 | -0.8 | 20.0 |
| PCB-1268 Peak 4 | Lin1 | | 0.0306 | | 0.490 | 0.500 | -2.0 | 20.0 |
| PCB-1268 Peak 5 | Lin1 | | 0.2305 | | 0.468 | 0.500 | -6.4 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206283/7  Calibration Date: 11/11/2015 12:06

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 19:46

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 21:08

Lab File ID: P4000007.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1221 Peak 1 | Lin1 | | 0.0071 | | 0.575 | 0.500 | 15.0 | 20.0 |
| PCB-1221 Peak 2 | Lin1 | | 0.0048 | | 0.590 | 0.500 | 18.0 | 20.0 |
| PCB-1221 Peak 3 | Lin1 | | 0.0173 | | 0.565 | 0.500 | 13.1 | 20.0 |
| PCB-1254 Peak 1 | Lin1 | | 0.0240 | | 0.556 | 0.500 | 11.2 | 20.0 |
| PCB-1254 Peak 2 | Lin1 | | 0.0379 | | 0.530 | 0.500 | 6.1 | 20.0 |
| PCB-1254 Peak 3 | Lin1 | | 0.0280 | | 0.537 | 0.500 | 7.4 | 20.0 |
| PCB-1254 Peak 4 | Lin1 | | 0.0225 | | 0.526 | 0.500 | 5.1 | 20.0 |
| PCB-1254 Peak 5 | Lin1 | | 0.0372 | | 0.528 | 0.500 | 5.6 | 20.0 |

FORM VII
PCBS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVIS 240-206283/28  Calibration Date: 11/11/2015 17:50

Instrument ID: A2HP4  Calib Start Date: 11/04/2015 21:24

GC Column: CLP-2  ID: 0.53(mm)  Calib End Date: 11/04/2015 22:46

Lab File ID: P4000028.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE RRF | RRF | MIN RRF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| PCB-1016 Peak 1 | Lin1 | | 0.0144 | | 0.541 | 0.500 | 8.2 | 20.0 |
| PCB-1016 Peak 2 | Lin1 | | 0.0279 | | 0.540 | 0.500 | 7.9 | 20.0 |
| PCB-1016 Peak 3 | Lin1 | | 0.0609 | | 0.517 | 0.500 | 3.4 | 20.0 |
| PCB-1016 Peak 4 | Lin1 | | 0.0243 | | 0.528 | 0.500 | 5.6 | 20.0 |
| PCB-1016 Peak 5 | Lin1 | | 0.0182 | | 0.519 | 0.500 | 3.9 | 20.0 |
| PCB-1260 Peak 1 | Lin1 | | 0.0327 | | 0.493 | 0.500 | -1.4 | 20.0 |
| PCB-1260 Peak 2 | Lin1 | | 0.0419 | | 0.500 | 0.500 | 0.0 | 20.0 |
| PCB-1260 Peak 3 | Lin1 | | 0.0623 | | 0.494 | 0.500 | -1.2 | 20.0 |
| PCB-1260 Peak 4 | Lin1 | | 0.0728 | | 0.482 | 0.500 | -3.5 | 20.0 |
| PCB-1260 Peak 5 | Lin1 | | 0.0501 | | 0.486 | 0.500 | -2.7 | 20.0 |
| Tetrachloro-m-xylene | Lin1 | | 0.6986 | | 0.0264 | 0.0250 | 5.8 | 20.0 |
| DCB Decachlorobiphenyl | Lin1 | | 0.5154 | | 0.0219 | 0.0250 | -12.3 | 20.0 |

# PCBS BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205809  Batch Start Date: 11/09/15 11:00  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/10/15 03:00

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | ex10PPMSPK 00018 | ex2/.2SURRW 00062 | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-E-1 | F-VB-UST-1-14 | 3540C, 8082A | T | 29.54 g | 10 mL | | 1 mL | | |
| 240-57597-D-2 | F-VB-UST-2-14 | 3540C, 8082A | T | 30.07 g | 10 mL | | 1 mL | | |
| 240-57597-E-3 | F-VB-UST-3-14 | 3540C, 8082A | T | 30.14 g | 10 mL | | 1 mL | | |
| 240-57597-E-4 | F-VB-UST-4-14 | 3540C, 8082A | T | 29.54 g | 10 mL | | 1 mL | | |
| 240-57597-E-5 | F-VB-UST-5-14 | 3540C, 8082A | T | 30.48 g | 10 mL | | 1 mL | | |
| 240-57597-E-6 | F-VB-UST-6-14 | 3540C, 8082A | T | 29.87 g | 10 mL | | 1 mL | | |
| 240-57597-E-7 | F-VB-UST-7-14 | 3540C, 8082A | T | 30.04 g | 10 mL | | 1 mL | | |
| 240-57597-D-8 | F-VB-UST-8-14 | 3540C, 8082A | T | 29.74 g | 10 mL | | 1 mL | | |
| 240-57597-E-9 | F-VB-UST-9-14 | 3540C, 8082A | T | 30.07 g | 10 mL | | 1 mL | | |
| 240-57597-E-10 | F-VB-UST-10-14 | 3540C, 8082A | T | 29.97 g | 10 mL | | 1 mL | | |
| 240-57597-D-11 | F-VB-UST-11-14 | 3540C, 8082A | T | 30.28 g | 10 mL | | 1 mL | | |
| 240-57597-D-12 | F-VB-UST-12-14 | 3540C, 8082A | T | 30.29 g | 10 mL | | 1 mL | | |
| 240-57597-D-13 | F-VB-UST-13-14 | 3540C, 8082A | T | 30.41 g | 10 mL | | 1 mL | | |
| 240-57597-E-14 | F-VB-UST-14-14 | 3540C, 8082A | T | 30.19 g | 10 mL | | 1 mL | | |
| 240-57597-E-15 | F-VB-UST-DUP | 3540C, 8082A | T | 30.34 g | 10 mL | | 1 mL | | |
| 240-57597-E-15 MS | F-VB-UST-DUP | 3540C, 8082A | T | 29.65 g | 10 mL | 1 mL | 1 mL | | |
| 240-57597-E-15 MSD | F-VB-UST-DUP | 3540C, 8082A | T | 29.78 g | 10 mL | 1 mL | 1 mL | | |
| MB 240-205809/21 | | 3540C, 8082A | | 30 g | 10 mL | | 1 mL | | |
| LCS 240-205809/22 | | 3540C, 8082A | | 30 g | 10 mL | 1 mL | 1 mL | | |

| Batch Notes | |
|---|---|
| Acid used for Clean Up Reagent | 2226515 |
| Balance ID | 1339912 |
| Na2SO4 Lot Number | 2204510 |
| Prep Solvent Lot # | 2247000 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

FORM II
GC VOA SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): RTX-VRX ID: 0.45(mm)

| Client Sample ID | Lab Sample ID | TFT1 # |
|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 78 |
| F-VB-UST-2-14 | 240-57597-2 | 79 |
| F-VB-UST-3-14 | 240-57597-3 | 77 |
| F-VB-UST-4-14 | 240-57597-4 | 71 |
| F-VB-UST-5-14 | 240-57597-5 | 71 |
| F-VB-UST-6-14 | 240-57597-6 | 76 |
| F-VB-UST-7-14 | 240-57597-7 | 78 |
| F-VB-UST-8-14 | 240-57597-8 | 73 |
| F-VB-UST-9-14 | 240-57597-9 | 78 |
| F-VB-UST-10-14 | 240-57597-10 | 79 |
| F-VB-UST-11-14 | 240-57597-11 | 79 |
| F-VB-UST-12-14 | 240-57597-12 | 78 |
| F-VB-UST-13-14 | 240-57597-13 | 75 |
| F-VB-UST-14-14 | 240-57597-14 | 80 |
| F-VB-UST-DUP | 240-57597-15 | 74 |
| | MB 240-206227/1-A | 73 |
| | MB 240-206444/1-A | 72 |
| | LCS 240-206227/2-A | 95 |
| | LCS 240-206444/2-A | 100 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 99 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 40 |

| | QC LIMITS |
|---|---|
| TFT = Trifluorotoluene (Surr) | 40-139 |

# Column to be used to flag recovery values

FORM II 8015C

FORM IV
GC VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: MB 240-206227/1-A

Matrix: Solid  Date Extracted: 11/11/2015 06:58

Lab File ID:(1) A0111107.D  Lab File ID:(2)

Date Analyzed:(1) 11/11/2015 09:56  Date Analyzed:(2)

Instrument ID:(1) AFID  Instrument ID:(2)

GC Column:(1) RTX-VRX  ID: 0.45(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-206227/2-A | 11/11/2015 10:31 | |
| F-VB-UST-1-14 | 240-57597-1 | 11/11/2015 11:08 | |
| F-VB-UST-2-14 | 240-57597-2 | 11/11/2015 11:45 | |
| F-VB-UST-3-14 | 240-57597-3 | 11/11/2015 12:23 | |
| F-VB-UST-4-14 | 240-57597-4 | 11/11/2015 13:00 | |
| F-VB-UST-6-14 | 240-57597-6 | 11/11/2015 14:13 | |
| F-VB-UST-7-14 | 240-57597-7 | 11/11/2015 14:50 | |
| F-VB-UST-8-14 | 240-57597-8 | 11/11/2015 15:27 | |
| F-VB-UST-9-14 | 240-57597-9 | 11/11/2015 17:19 | |
| F-VB-UST-10-14 | 240-57597-10 | 11/11/2015 17:57 | |
| F-VB-UST-11-14 | 240-57597-11 | 11/11/2015 18:34 | |
| F-VB-UST-12-14 | 240-57597-12 | 11/11/2015 19:12 | |
| F-VB-UST-13-14 | 240-57597-13 | 11/11/2015 19:48 | |
| F-VB-UST-14-14 | 240-57597-14 | 11/11/2015 20:25 | |
| F-VB-UST-DUP | 240-57597-15 | 11/11/2015 21:03 | |
| F-VB-UST-DUP MS | 240-57597-15 MS | 11/11/2015 21:40 | |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 11/11/2015 22:17 | |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-206227/1-A

Matrix: Solid  Lab File ID: A0111107.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/11/2015 09:56

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206211  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 46 | U | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 73 | | 40-139 |

FORM III
GC VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111108.D

Lab ID: LCS 240-206227/2-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 786 | 98 | 74-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-206227/2-A

Matrix: Solid  Lab File ID: A0111108.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/11/2015 10:31

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206211  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 786 | | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 95 | | 40-139 |

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\AFID\20151111-49133.b\A0111108.D
Lims ID: LCS 240-206227/2-A
Client ID:
Sample Type: LCS
Inject. Date: 11-Nov-2015 10:31:41 ALS Bottle#: 8 Worklist Smp#: 8
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Sample Info: 240-0049133-008
Operator ID: Instrument ID: AFID

Method: \\ChromNA\Canton\ChromData\AFID\20151111-49133.b\A80218015.m
Limit Group: GCV 8015C GRO ICAL
Last Update: 12-Nov-2015 06:28:18 Calib Date: 24-Aug-2015 17:07:27
Integrator: Falcon
Quant Method: External Standard Quant By: Initial Calibration
Last ICal File: \\ChromNA\Canton\ChromData\AFID\20150824-46596.b\A0082112.D

Column 1 : Det: GC FID1A
Process Host: XAWRK032

| Compound | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Response | Cal Amt ug/l | OnCol Amt ug/l | Flags |
|---|---|---|---|---|---|---|---|
| $ 16 Trifluorotoluene (Surr) | 10.363 | 10.366 | -0.003 | 4201616 | 10.0 | 9.46 | |
| A 5 C6-C10 | 11.535 | (5.939-17.178) | | 318414423 | 800.0 | 786.4 | |
| A 1 C6-C12 | 12.801 | (5.939-19.699) | | 363011179 | 800.0 | 806.4 | |

## Reagents:

VGGASCHK_00070 Amount Added: 40.00 Units: uL
VGGASURROGATE_00034 Amount Added: 1.00 Units: uL Run Reagent

FORM III
GC VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111126.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (ug/Kg) | SAMPLE CONCENTRATION (ug/Kg) | MS CONCENTRATION (ug/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 928 | 54 U | 856 | 92 | 10-136 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
GC VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Matrix: Solid  Level: Low  Lab File ID: A0111127.D

Lab ID: 240-57597-15 MSD  Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (ug/Kg) | MSD CONCENTRATION (ug/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 926 | 206 | 22 | 122 | 40 | 10-136 | F2 |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MS | Lab Sample ID: 240-57597-15 MS

Matrix: Solid | Lab File ID: A0111126.D

Analysis Method: 8015C | Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.04(g) | Date Analyzed: 11/11/2015 21:40

Soil Aliquot Vol: | Dilution Factor: 1

Soil Extract Vol.: | GC Column: RTX-VRX ID: 0.45(mm)

% Moisture: 14.5 | Level: (low/med) Low

Analysis Batch No.: 206211 | Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 856 | | 120 | 53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 99 | | 40-139 |

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\AFID\20151111-49133.b\A0111126.D
Lims ID: 240-57597-C-15-B MS
Client ID:
Sample Type: MS
Inject. Date: 11-Nov-2015 21:40:33 ALS Bottle#: 26 Worklist Smp#: 26
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Sample Info: 240-0049133-026
Operator ID: Instrument ID: AFID

Method: \\ChromNA\Canton\ChromData\AFID\20151111-49133.b\A80218015.m
Limit Group: GCV 8015C GRO ICAL
Last Update: 12-Nov-2015 06:28:29 Calib Date: 24-Aug-2015 17:07:27
Integrator: Falcon
Quant Method: External Standard Quant By: Initial Calibration
Last ICal File: \\ChromNA\Canton\ChromData\AFID\20150824-46596.b\A0082112.D

Column 1 : Det: GC FID1A
Process Host: XAWRK032

| Compound | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Response | Cal Amt ug/l | OnCol Amt ug/l | Flags |
|---|---|---|---|---|---|---|---|
| 6 2-Methylpentane | 6.088 | 6.101 | -0.013 | 14862522 | | NC | |
| 12 Methyl tert-butyl ether | 6.453 | 6.421 | 0.032 | 7414201 | | NC | |
| $ 16 Trifluorotoluene (Surr) | 10.368 | 10.374 | -0.006 | 4390464 | 10.0 | 9.88 | |
| A 5 C6-C10 | 11.559 | (5.939-17.178) | | 298900015 | 800.0 | 738.2 | |
| A 1 C6-C12 | 12.819 | (5.939-19.699) | | 335788384 | 800.0 | 745.9 | |
| A 11 GRO | 13.152 | (6.265-20.038) | | 350425266 | | NC | |
| 17 1,2,4-Trimethylbenzene | 17.011 | 17.010 | 0.001 | 12491028 | | NC | |
| 14 Dodecane | 19.505 | 19.498 | 0.007 | 706013 | | NC | |
| 9 Naphthalene | 19.843 | 19.888 | -0.045 | 1526118 | | NC | |

## QC Flag Legend

Processing Flags
NC - Not Calibrated

## Reagents:

VGGASCHK_00070 Amount Added: 40.00 Units: uL
VGGASURROGATE_00034 Amount Added: 1.00 Units: uL Run Reagent

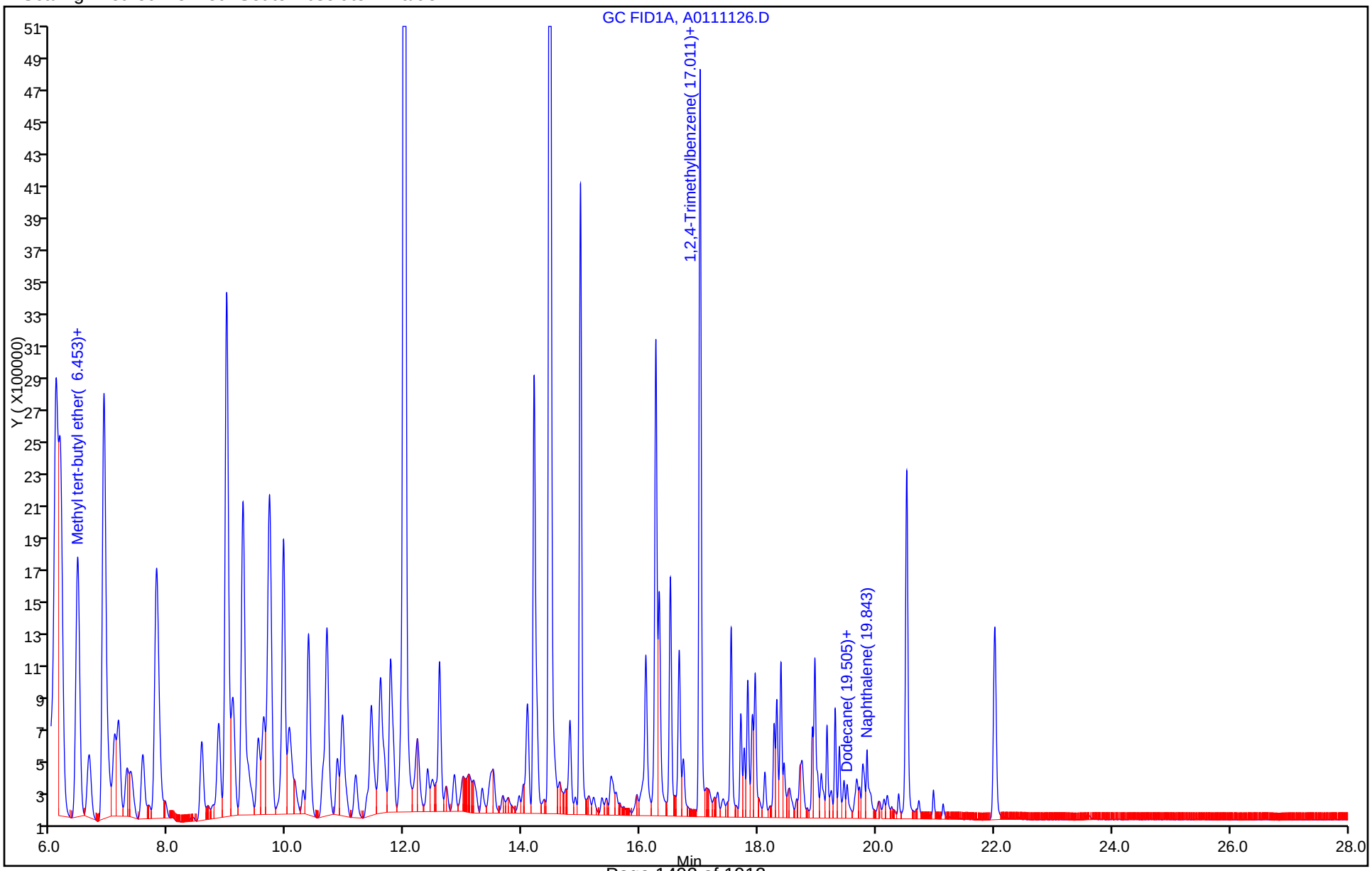
Report Date: 12-Nov-2015 06:28:35 Chrom Revision: 2.2 08-Oct-2015 07:17:48

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\AFID\20151111-49133.b\A0111126.D

| | | | | | |
|---|---|---|---|---|---|
| Injection Date: | 11-Nov-2015 21:40:33 | Instrument ID: | AFID | Operator ID: | |
| Lims ID: | 240-57597-C-15-B MS | | | Worklist Smp#: | 26 |
| Client ID: | | | | | |
| Purge Vol: | 5.000 mL | Dil. Factor: | 1.0000 | ALS Bottle#: | 26 |
| Method: | A80218015 | Limit Group: | GCV 8015C GRO ICAL | | |

Y Scaling: Method Defined: Set to Absolute Y Value



FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID: F-VB-UST-DUP MSD  Lab Sample ID: 240-57597-15 MSD

Matrix: Solid  Lab File ID: A0111127.D

Analysis Method: 8015C  Date Collected: 11/05/2015 00:00

Sample wt/vol: 5.05(g)  Date Analyzed: 11/11/2015 22:17

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX  ID: 0.45(mm)

% Moisture: 14.5  Level: (low/med) Low

Analysis Batch No.: 206211  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 206 | | 120 | 53 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 40 | | 40-139 |

MSD %R = 22%

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\AFID\20151111-49133.b\A0111127.D
Lims ID: 240-57597-C-15-C MSD
Client ID:
Sample Type: MSD
Inject. Date: 11-Nov-2015 22:17:52 ALS Bottle#: 27 Worklist Smp#: 27
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Sample Info: 240-0049133-027
Operator ID: Instrument ID: AFID

Method: \\ChromNA\Canton\ChromData\AFID\20151111-49133.b\A80218015.m
Limit Group: GCV 8015C GRO ICAL
Last Update: 12-Nov-2015 06:28:29 Calib Date: 24-Aug-2015 17:07:27
Integrator: Falcon
Quant Method: External Standard Quant By: Initial Calibration
Last ICal File: \\ChromNA\Canton\ChromData\AFID\20150824-46596.b\A0082112.D

Column 1 : Det: GC FID1A
Process Host: XAWRK032

| Compound | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Response | Cal Amt ug/l | OnCol Amt ug/l | Flags |
|---|---|---|---|---|---|---|---|
| 6 2-Methylpentane | 6.139 | 6.101 | 0.038 | 7317419 | | NC | |
| 12 Methyl tert-butyl ether | 6.442 | 6.421 | 0.021 | 1636776 | | NC | |
| $ 16 Trifluorotoluene (Surr) | 10.367 | 10.374 | -0.007 | 1789897 | 10.0 | 4.03 | |
| A 5 C6-C10 | 11.559 | (5.939-17.178) | | 72198879 | 800.0 | 178.3 | |
| A 1 C6-C12 | 12.819 | (5.939-19.699) | | 81609255 | 800.0 | 181.3 | |
| A 11 GRO | 13.152 | (6.265-20.038) | | 99483778 | | NC | |
| 17 1,2,4-Trimethylbenzene | 17.014 | 17.010 | 0.004 | 3602706 | | NC | |
| 14 Dodecane | 19.509 | 19.498 | 0.011 | 143429 | | NC | |
| 9 Naphthalene | 19.909 | 19.888 | 0.021 | 76360 | | NC | |

## QC Flag Legend

Processing Flags
NC - Not Calibrated

## Reagents:

VGGASCHK_00070 Amount Added: 40.00 Units: uL
VGGASURROGATE_00034 Amount Added: 1.00 Units: uL Run Reagent

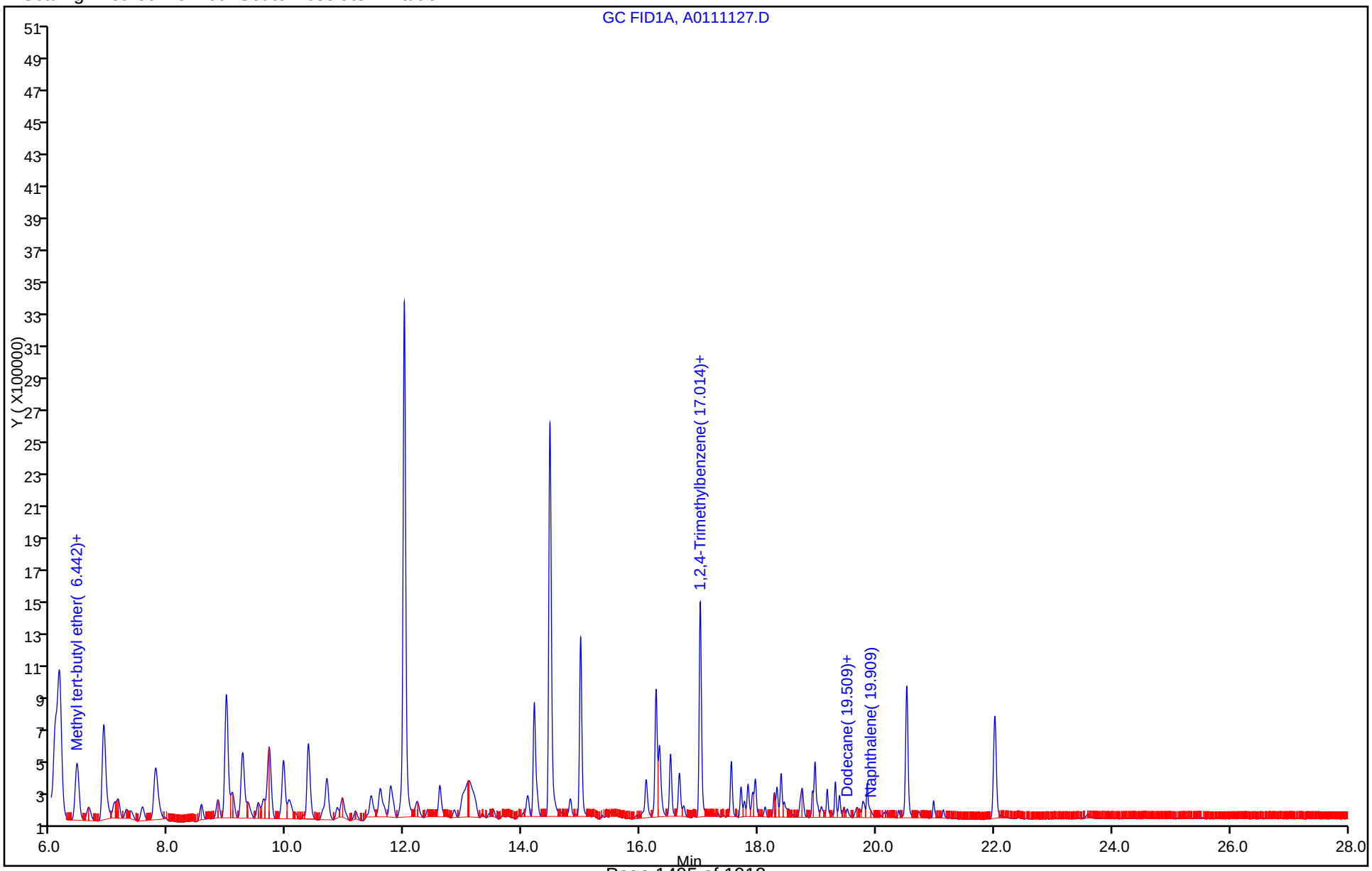
Report Date: 12-Nov-2015 06:28:36 Chrom Revision: 2.2 08-Oct-2015 07:17:48

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\AFID\20151111-49133.b\A0111127.D

| | | | | | |
|---|---|---|---|---|---|
| Injection Date: | 11-Nov-2015 22:17:52 | Instrument ID: | AFID | Operator ID: | |
| Lims ID: | 240-57597-C-15-C MSD | | | Worklist Smp#: | 27 |
| Client ID: | | | | | |
| Purge Vol: | 5.000 mL | Dil. Factor: | 1.0000 | ALS Bottle#: | 27 |
| Method: | A80218015 | Limit Group: | GCV 8015C GRO ICAL | | |

Y Scaling: Method Defined: Set to Absolute Y Value



FORM IV
GC VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: MB 240-206444/1-A

Matrix: Solid  Date Extracted: 11/12/2015 05:59

Lab File ID:(1) A0111205.D  Lab File ID:(2)

Date Analyzed:(1) 11/12/2015 08:03  Date Analyzed:(2)

Instrument ID:(1) AFID  Instrument ID:(2)

GC Column:(1) RTX-VRX  ID: 0.45(mm)  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-206444/2-A | 11/12/2015 08:40 | |
| F-VB-UST-5-14 | 240-57597-5 | 11/12/2015 09:17 | |

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: MB 240-206444/1-A

Matrix: Solid  Lab File ID: A0111205.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/12/2015 08:03

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206404  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 46 | U | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 72 | | 40-139 |

FORM III
GC VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: A0111206.D

Lab ID: LCS 240-206444/2-A Client ID:

| COMPOUND | SPIKE ADDED (ug/Kg) | LCS CONCENTRATION (ug/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | 800 | 787 | 98 | 74-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM I
GC VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Client Sample ID:  Lab Sample ID: LCS 240-206444/2-A

Matrix: Solid  Lab File ID: A0111206.D

Analysis Method: 8015C  Date Collected:

Sample wt/vol: 5(g)  Date Analyzed: 11/12/2015 08:40

Soil Aliquot Vol:  Dilution Factor: 1

Soil Extract Vol.:  GC Column: RTX-VRX ID: 0.45(mm)

% Moisture:  Level: (low/med) Low

Analysis Batch No.: 206404  Units: ug/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| 8006-61-9 | Gasoline Range Organics [C6 - C10] | 787 | | 100 | 46 |

| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 98-08-8 | Trifluorotoluene (Surr) | 100 | | 40-139 |

TestAmerica Canton
Target Compound Quantitation Report

Data File: \\ChromNA\Canton\ChromData\AFID\20151112-49174.b\A0111206.D
Lims ID: LCS 240-206444/2-A
Client ID:
Sample Type: LCS
Inject. Date: 12-Nov-2015 08:40:00 ALS Bottle#: 6 Worklist Smp#: 6
Purge Vol: 5.000 mL Dil. Factor: 1.0000
Sample Info:
Operator ID: Instrument ID: AFID

Method: \\ChromNA\Canton\ChromData\AFID\20151112-49174.b\A80218015.m
Limit Group: GCV 8015C GRO ICAL
Last Update: 12-Nov-2015 09:30:49 Calib Date: 24-Aug-2015 17:07:27
Integrator: Falcon
Quant Method: External Standard Quant By: Initial Calibration
Last ICal File: \\ChromNA\Canton\ChromData\AFID\20150824-46596.b\A0082112.D

Column 1 : Det: GC FID1A
Process Host: XAWRK031

| Compound | RT (min.) | Exp RT (min.) | Dlt RT (min.) | Response | Cal Amt ug/l | OnCol Amt ug/l | Flags |
|---|---|---|---|---|---|---|---|
| $ 16 Trifluorotoluene (Surr) | 10.372 | 10.368 | 0.004 | 4430972 | 10.0 | 9.97 | |
| A 5 C6-C10 | 11.532 | (5.929-17.135) | | 318752592 | 800.0 | 787.2 | |
| A 1 C6-C12 | 12.792 | (5.929-19.655) | | 361056203 | 800.0 | 802.1 | |

## Reagents:

VGGASCHK_00070 Amount Added: 40.00 Units: uL
VGGASURROGATE_00034 Amount Added: 1.00 Units: uL Run Reagent

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: STD4 240-194508/10  Date Analyzed: 08/24/2015 15:53

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): A0082110.D  Heated Purge: (Y/N) N

Calibration ID: 30361

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| INITIAL CALIBRATION SURROGATE | | | | 10.34 | | |
| UPPER LIMIT | | | | 10.44 | | |
| LOWER LIMIT | | | | 10.24 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| STD4 240-194508/10 ICRT | | 08/24/2015 15:53 | A0082110.D | 10.34 | | |
| ICV 240-194508/13 | | 08/24/2015 17:45 | A0082113.D | 10.34 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVRT 240-206211/4  Date Analyzed: 11/11/2015 08:08

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): A0111104.D  Heated Purge: (Y/N) N

Calibration ID: 30361

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT RT # | | |
|---|---|---|---|---|---|---|
| CONTINUING CALIBRATION SURROGATE | | | | 10.36 | | |
| UPPER LIMIT | | | | 10.46 | | |
| LOWER LIMIT | | | | 10.26 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCVRT 240-206211/4 | | 11/11/2015 08:08 | A0111104.D | 10.36 | | |
| CCV 240-206211/6 | | 11/11/2015 09:20 | A0111106.D | 10.37 | | |
| MB 240-206227/1-A | | 11/11/2015 09:56 | A0111107.D | 10.36 | | |
| LCS 240-206227/2-A | | 11/11/2015 10:31 | A0111108.D | 10.36 | | |
| 240-57597-1 | F-VB-UST-1-14 | 11/11/2015 11:08 | A0111109.D | 10.37 | | |
| 240-57597-2 | F-VB-UST-2-14 | 11/11/2015 11:45 | A0111110.D | 10.37 | | |
| 240-57597-3 | F-VB-UST-3-14 | 11/11/2015 12:23 | A0111111.D | 10.37 | | |
| 240-57597-4 | F-VB-UST-4-14 | 11/11/2015 13:00 | A0111112.D | 10.37 | | |
| 240-57597-6 | F-VB-UST-6-14 | 11/11/2015 14:13 | A0111114.D | 10.37 | | |
| 240-57597-7 | F-VB-UST-7-14 | 11/11/2015 14:50 | A0111115.D | 10.37 | | |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 15:27 | A0111116.D | 10.37 | | |
| CCV 240-206211/18 | | 11/11/2015 16:42 | A0111118.D | 10.37 | | |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 17:19 | A0111119.D | 10.37 | | |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 17:57 | A0111120.D | 10.37 | | |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 18:34 | A0111121.D | 10.37 | | |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 19:12 | A0111122.D | 10.37 | | |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 19:48 | A0111123.D | 10.37 | | |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 20:25 | A0111124.D | 10.37 | | |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 21:03 | A0111125.D | 10.37 | | |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 21:40 | A0111126.D | 10.37 | | |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 22:17 | A0111127.D | 10.37 | | |
| CCV 240-206211/28 | | 11/11/2015 22:54 | A0111128.D | 10.37 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
GC VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCVRT 240-206404/4  Date Analyzed: 11/12/2015 07:26

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)

Lab File ID (Standard): A0111204.D  Heated Purge: (Y/N) N

Calibration ID: 30361

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | TFT | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 10.37 | | |
| UPPER LIMIT | | | | 10.47 | | |
| LOWER LIMIT | | | | 10.27 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCVRT 240-206404/4 | | 11/12/2015 07:26 | A0111204.D | 10.37 | | |
| MB 240-206444/1-A | | 11/12/2015 08:03 | A0111205.D | 10.37 | | |
| LCS 240-206444/2-A | | 11/12/2015 08:40 | A0111206.D | 10.37 | | |
| 240-57597-5 | F-VB-UST-5-14 | 11/12/2015 09:17 | A0111207.D | 10.37 | | |
| CCV 240-206404/8 | | 11/12/2015 09:54 | A0111208.D | 10.37 | | |

TFT = Trifluorotoluene (Surr)

TFT RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Instrument ID: AFID  Start Date: 08/24/2015 13:59

Analysis Batch Number: 194508  End Date: 08/24/2015 17:45

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-194508/7 IC | | 08/24/2015 13:59 | 1 | A0082107.D | RTX-VRX 0.45(mm) |
| STD2 240-194508/8 IC | | 08/24/2015 14:37 | 1 | A0082108.D | RTX-VRX 0.45(mm) |
| STD3 240-194508/9 IC | | 08/24/2015 15:14 | 1 | A0082109.D | RTX-VRX 0.45(mm) |
| STD4 240-194508/10 ICRT | | 08/24/2015 15:53 | 1 | A0082110.D | RTX-VRX 0.45(mm) |
| STD5 240-194508/11 IC | | 08/24/2015 16:31 | 1 | A0082111.D | RTX-VRX 0.45(mm) |
| STD6 240-194508/12 IC | | 08/24/2015 17:07 | 1 | A0082112.D | RTX-VRX 0.45(mm) |
| ICV 240-194508/13 | | 08/24/2015 17:45 | 1 | A0082113.D | RTX-VRX 0.45(mm) |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: AFID Start Date: 11/11/2015 08:08

Analysis Batch Number: 206211 End Date: 11/11/2015 22:54

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCVRT 240-206211/4 | | 11/11/2015 08:08 | 1 | A0111104.D | RTX-VRX 0.45(mm) |
| CCV 240-206211/6 | | 11/11/2015 09:20 | 1 | A0111106.D | RTX-VRX 0.45(mm) |
| MB 240-206227/1-A | | 11/11/2015 09:56 | 1 | A0111107.D | RTX-VRX 0.45(mm) |
| LCS 240-206227/2-A | | 11/11/2015 10:31 | 1 | A0111108.D | RTX-VRX 0.45(mm) |
| 240-57597-1 | F-VB-UST-1-14 | 11/11/2015 11:08 | 1 | A0111109.D | RTX-VRX 0.45(mm) |
| 240-57597-2 | F-VB-UST-2-14 | 11/11/2015 11:45 | 1 | A0111110.D | RTX-VRX 0.45(mm) |
| 240-57597-3 | F-VB-UST-3-14 | 11/11/2015 12:23 | 1 | A0111111.D | RTX-VRX 0.45(mm) |
| 240-57597-4 | F-VB-UST-4-14 | 11/11/2015 13:00 | 1 | A0111112.D | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/11/2015 13:36 | 1 | | RTX-VRX 0.45(mm) |
| 240-57597-6 | F-VB-UST-6-14 | 11/11/2015 14:13 | 1 | A0111114.D | RTX-VRX 0.45(mm) |
| 240-57597-7 | F-VB-UST-7-14 | 11/11/2015 14:50 | 1 | A0111115.D | RTX-VRX 0.45(mm) |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 15:27 | 1 | A0111116.D | RTX-VRX 0.45(mm) |
| CCV 240-206211/18 | | 11/11/2015 16:42 | 1 | A0111118.D | RTX-VRX 0.45(mm) |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 17:19 | 1 | A0111119.D | RTX-VRX 0.45(mm) |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 17:57 | 1 | A0111120.D | RTX-VRX 0.45(mm) |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 18:34 | 1 | A0111121.D | RTX-VRX 0.45(mm) |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 19:12 | 1 | A0111122.D | RTX-VRX 0.45(mm) |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 19:48 | 1 | A0111123.D | RTX-VRX 0.45(mm) |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 20:25 | 1 | A0111124.D | RTX-VRX 0.45(mm) |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 21:03 | 1 | A0111125.D | RTX-VRX 0.45(mm) |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 21:40 | 1 | A0111126.D | RTX-VRX 0.45(mm) |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 22:17 | 1 | A0111127.D | RTX-VRX 0.45(mm) |
| CCV 240-206211/28 | | 11/11/2015 22:54 | 1 | A0111128.D | RTX-VRX 0.45(mm) |

GC VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: AFID Start Date: 11/12/2015 07:26

Analysis Batch Number: 206404 End Date: 11/13/2015 01:22

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCVRT 240-206404/4 | | 11/12/2015 07:26 | 1 | A0111204.D | RTX-VRX 0.45(mm) |
| MB 240-206444/1-A | | 11/12/2015 08:03 | 1 | A0111205.D | RTX-VRX 0.45(mm) |
| LCS 240-206444/2-A | | 11/12/2015 08:40 | 1 | A0111206.D | RTX-VRX 0.45(mm) |
| 240-57597-5 | F-VB-UST-5-14 | 11/12/2015 09:17 | 1 | A0111207.D | RTX-VRX 0.45(mm) |
| CCV 240-206404/8 | | 11/12/2015 09:54 | 1 | A0111208.D | RTX-VRX 0.45(mm) |
| CCV 240-206404/16 | | 11/12/2015 14:54 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/12/2015 15:31 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/12/2015 16:09 | 1 | | RTX-VRX 0.45(mm) |
| CCV 240-206404/22 | | 11/12/2015 18:38 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/12/2015 19:51 | 1 | | RTX-VRX 0.45(mm) |
| ZZZZZ | | 11/12/2015 20:28 | 1 | | RTX-VRX 0.45(mm) |
| CCV 240-206404/33 | | 11/13/2015 01:22 | 1 | | RTX-VRX 0.45(mm) |

8015C

FORM VI
GC VOA BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton  Job No.: 240-57597-1  Analy Batch No.: 194508

SDG No.:

Instrument ID: AFID  GC Column: RTX-VRX  ID: 0.45(mm)  Heated Purge: (Y/N) N

Calibration Start Date: 08/24/2015 13:59  Calibration End Date: 08/24/2015 17:07  Calibration ID: 30361

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-194508/7 | A0082107.D |
| Level 2 | STD2 240-194508/8 | A0082108.D |
| Level 3 | STD3 240-194508/9 | A0082109.D |
| Level 4 | STD4 240-194508/10 | A0082110.D |
| Level 5 | STD5 240-194508/11 | A0082111.D |
| Level 6 | STD6 240-194508/12 | A0082112.D |

| ANALYTE | CF | | | | CURVE TYPE | COEFFICIENT | | | # | MIN CF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1<br>LVL 5 | LVL 2<br>LVL 6 | LVL 3 | LVL 4 | | B | M1 | M2 | | | | | | | | |
| Gasoline Range Organics [C6 - C10] | 281051<br>441376 | 404289<br>439409 | 445616 | 417800 | Ave | | 404923.648 | | | | 15.5 | | 20.0 | | | |
| C6-C12 | 344884<br>484706 | 442549<br>478304 | 494744 | 455788 | Ave | | 450162.318 | | | | 12.2 | | 20.0 | | | |
| Trifluorotoluene (Surr) | 452208<br>466474 | 509634<br>371094 | 436510 | 429984 | Ave | | 444317.252 | | | | 10.3 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-194508/13  Calibration Date: 08/24/2015 17:45

Instrument ID: AFID  Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX  ID: 0.45(mm)  Calib End Date: 08/24/2015 17:07

Lab File ID: A0082113.D  Conc. Units: ug/L  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 425225 | | 840 | 800 | 5.0 | 20.0 |
| C6-C12 | Ave | 450162 | 466294 | | 829 | 800 | 3.6 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 487053 | | 11.0 | 10.0 | 9.6 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVRT 240-206211/4 Calibration Date: 11/11/2015 08:08

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111104.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 420445 | | 831 | 800 | 3.8 | 20.0 |
| C6-C12 | Ave | 450162 | 459766 | | 817 | 800 | 2.1 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 408705 | | 9.20 | 10.0 | -8.0 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206211/6  Calibration Date: 11/11/2015 09:20

Instrument ID: AFID  Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX  ID: 0.45(mm)  Calib End Date: 08/24/2015 17:07

Lab File ID: A0111106.D  Conc. Units: ug/L  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 428576 | | 847 | 800 | 5.8 | 20.0 |
| C6-C12 | Ave | 450162 | 466289 | | 829 | 800 | 3.6 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 402849 | | 9.07 | 10.0 | -9.3 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206211/18 Calibration Date: 11/11/2015 16:42

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111118.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 421224 | | 832 | 800 | 4.0 | 20.0 |
| C6-C12 | Ave | 450162 | 459626 | | 817 | 800 | 2.1 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 402080 | | 9.05 | 10.0 | -9.5 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206211/28 Calibration Date: 11/11/2015 22:54

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111128.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 417296 | | 824 | 800 | 3.1 | 20.0 |
| C6-C12 | Ave | 450162 | 454853 | | 808 | 800 | 1.0 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 401488 | | 9.04 | 10.0 | -9.6 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCVRT 240-206404/4 Calibration Date: 11/12/2015 07:26

Instrument ID: AFID Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX ID: 0.45(mm) Calib End Date: 08/24/2015 17:07

Lab File ID: A0111204.D Conc. Units: ug/L Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 427634 | | 845 | 800 | 5.6 | 20.0 |
| C6-C12 | Ave | 450162 | 465474 | | 827 | 800 | 3.4 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 405884 | | 9.14 | 10.0 | -8.6 | 20.0 |

FORM VII
GC VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206404/8  Calibration Date: 11/12/2015 09:54

Instrument ID: AFID  Calib Start Date: 08/24/2015 13:59

GC Column: RTX-VRX  ID: 0.45(mm)  Calib End Date: 08/24/2015 17:07

Lab File ID: A0111208.D  Conc. Units: ug/L  Heated Purge: (Y/N) N

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Gasoline Range Organics [C6 - C10] | Ave | 404924 | 428972 | | 848 | 800 | 5.9 | 20.0 |
| C6-C12 | Ave | 450162 | 465297 | | 827 | 800 | 3.4 | 20.0 |
| Trifluorotoluene (Surr) | Ave | 444317 | 420936 | | 9.47 | 10.0 | -5.3 | 20.0 |

# GC VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 206227  Batch Start Date: 11/11/15 06:58  Batch Analyst: Reed, Rachel T

Batch Method: 5030A  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| MB 240-206227/1 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| LCS 240-206227/2 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| 240-57597-C-1 | F-VB-UST-1-14 | 5030A, 8015C | T | 4.98 g | 5 mL | | | | |
| 240-57597-C-2 | F-VB-UST-2-14 | 5030A, 8015C | T | 4.98 g | 5 mL | | | | |
| 240-57597-C-3 | F-VB-UST-3-14 | 5030A, 8015C | T | 4.99 g | 5 mL | | | | |
| 240-57597-C-4 | F-VB-UST-4-14 | 5030A, 8015C | T | 4.99 g | 5 mL | | | | |
| 240-57597-C-6 | F-VB-UST-6-14 | 5030A, 8015C | T | 5.01 g | 5 mL | | | | |
| 240-57597-C-7 | F-VB-UST-7-14 | 5030A, 8015C | T | 4.97 g | 5 mL | | | | |
| 240-57597-C-8 | F-VB-UST-8-14 | 5030A, 8015C | T | 4.98 g | 5 mL | | | | |
| 240-57597-C-9 | F-VB-UST-9-14 | 5030A, 8015C | T | 5.00 g | 5 mL | | | | |
| 240-57597-C-10 | F-VB-UST-10-14 | 5030A, 8015C | T | 5.02 g | 5 mL | | | | |
| 240-57597-C-11 | F-VB-UST-11-14 | 5030A, 8015C | T | 5.03 g | 5 mL | | | | |
| 240-57597-C-12 | F-VB-UST-12-14 | 5030A, 8015C | T | 5.04 g | 5 mL | | | | |
| 240-57597-C-13 | F-VB-UST-13-14 | 5030A, 8015C | T | 4.99 g | 5 mL | | | | |
| 240-57597-C-14 | F-VB-UST-14-14 | 5030A, 8015C | T | 5.02 g | 5 mL | | | | |
| 240-57597-C-15 | F-VB-UST-DUP | 5030A, 8015C | T | 5.00 g | 5 mL | | | | |
| 240-57597-C-15 MS | F-VB-UST-DUP | 5030A, 8015C | T | 5.04 g | 5 mL | | | | |
| 240-57597-C-15 MSD | F-VB-UST-DUP | 5030A, 8015C | T | 5.05 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| Balance ID | B045 |
| Blank Soil Lot Number | 1757979 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# GC VOA BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 206444  Batch Start Date: 11/12/15 07:59  Batch Analyst: Reed, Rachel T

Batch Method: 5030A  Batch End Date:

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | | | | |
|---|---|---|---|---|---|---|---|---|---|
| MB 240-206444/1 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| LCS 240-206444/2 | | 5030A, 8015C | | 5 g | 5 mL | | | | |
| 240-57597-C-5 | F-VB-UST-5-14 | 5030A, 8015C | T | 4.98 g | 5 mL | | | | |

| Batch Notes | |
|---|---|
| Balance ID | B045 |
| Blank Soil Lot Number | 1757979 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

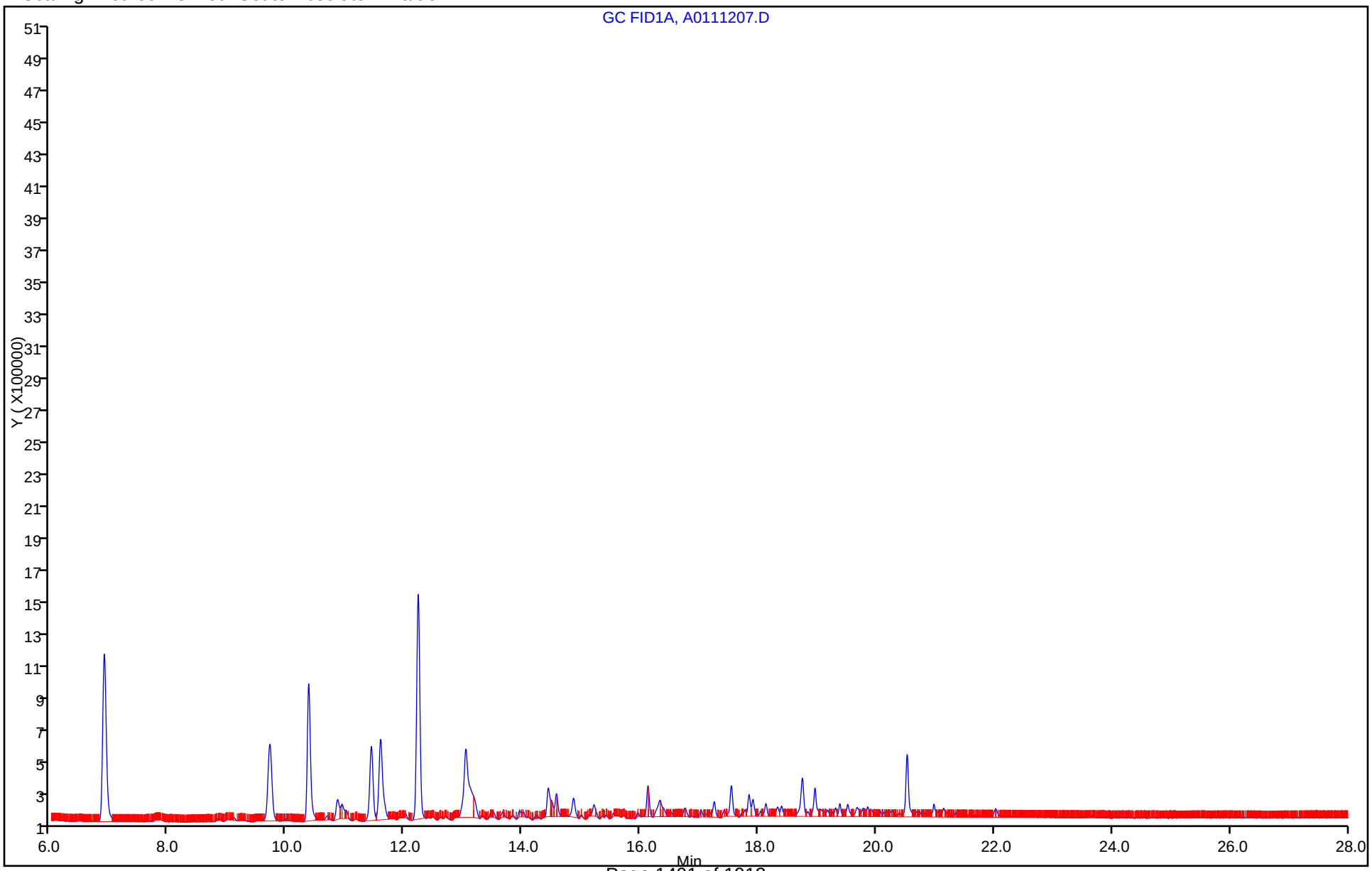
Report Date: 12-Nov-2015 09:30:55 Chrom Revision: 2.2 08-Oct-2015 07:17:48

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\AFID\20151112-49174.b\A0111207.D

| | | | | | |
|---|---|---|---|---|---|
| Injection Date: | 12-Nov-2015 09:17:06 | Instrument ID: | AFID | Operator ID: | |
| Lims ID: | 240-57597-C-5-B | Lab Sample ID: | 240-57597-5 | Worklist Smp#: | 7 |
| Client ID: | F-VB-UST-5-14 | | | | |
| Purge Vol: | 5.000 mL | Dil. Factor: | 1.0000 | ALS Bottle#: | 7 |
| Method: | A80218015 | Limit Group: | GCV 8015C GRO ICAL | | |

Y Scaling: Method Defined: Set to Absolute Y Value



FORM II
DIESEL RANGE ORGANICS SURROGATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low

GC Column (1): ID:

| Client Sample ID | Lab Sample ID | OTPH1 # |
|---|---|---|
| F-VB-UST-1-14 | 240-57597-1 | 62 |
| F-VB-UST-2-14 | 240-57597-2 | 62 |
| F-VB-UST-3-14 | 240-57597-3 | 59 |
| F-VB-UST-4-14 | 240-57597-4 | 60 |
| F-VB-UST-5-14 | 240-57597-5 | 56 |
| F-VB-UST-6-14 | 240-57597-6 | 60 |
| F-VB-UST-7-14 | 240-57597-7 | 65 |
| F-VB-UST-8-14 | 240-57597-8 | 65 |
| F-VB-UST-9-14 | 240-57597-9 | 57 |
| F-VB-UST-10-14 | 240-57597-10 | 64 |
| F-VB-UST-11-14 | 240-57597-11 | 64 |
| F-VB-UST-12-14 | 240-57597-12 | 73 |
| F-VB-UST-13-14 | 240-57597-13 | 68 |
| F-VB-UST-14-14 | 240-57597-14 | 61 |
| F-VB-UST-DUP | 240-57597-15 | 65 |
| | MB 240-205788/22-A | 56 |
| | LCS 240-205788/23-A | 67 |
| F-VB-UST-DUP MS | 240-57597-15 MS | 81 |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 73 |

| | QC LIMITS |
|---|---|
| OTPH = o-Terphenyl (Surr) | 40-160 |

# Column to be used to flag recovery values

FORM II 8015C

FORM IV
DIESEL RANGE ORGANICS METHOD BLANK SUMMARY

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: MB 240-205788/22-A

Matrix: Solid  Date Extracted: 11/09/2015 09:40

Lab File ID:(1) P5B11008.D  Lab File ID:(2)

Date Analyzed:(1) 11/10/2015 16:59  Date Analyzed:(2)

Instrument ID:(1) A2HP5R  Instrument ID:(2)

GC Column:(1)  ID:  GC Column:(2)  ID:

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

| CLIENT SAMPLE ID | LAB SAMPLE ID | DATE ANALYZED 1 | DATE ANALYZED 2 |
|---|---|---|---|
| | LCS 240-205788/23-A | 11/10/2015 17:31 | |
| F-VB-UST-1-14 | 240-57597-1 | 11/10/2015 20:07 | |
| F-VB-UST-2-14 | 240-57597-2 | 11/10/2015 20:38 | |
| F-VB-UST-3-14 | 240-57597-3 | 11/10/2015 21:10 | |
| F-VB-UST-4-14 | 240-57597-4 | 11/10/2015 21:41 | |
| F-VB-UST-5-14 | 240-57597-5 | 11/10/2015 22:43 | |
| F-VB-UST-6-14 | 240-57597-6 | 11/10/2015 23:14 | |
| F-VB-UST-7-14 | 240-57597-7 | 11/10/2015 23:46 | |
| F-VB-UST-8-14 | 240-57597-8 | 11/11/2015 00:17 | |
| F-VB-UST-9-14 | 240-57597-9 | 11/11/2015 00:48 | |
| F-VB-UST-10-14 | 240-57597-10 | 11/11/2015 01:19 | |
| F-VB-UST-11-14 | 240-57597-11 | 11/11/2015 01:50 | |
| F-VB-UST-12-14 | 240-57597-12 | 11/11/2015 19:03 | |
| F-VB-UST-13-14 | 240-57597-13 | 11/11/2015 19:34 | |
| F-VB-UST-14-14 | 240-57597-14 | 11/11/2015 20:06 | |
| F-VB-UST-DUP | 240-57597-15 | 11/11/2015 20:37 | |
| F-VB-UST-DUP MS | 240-57597-15 MS | 11/11/2015 21:08 | |
| F-VB-UST-DUP MSD | 240-57597-15 MSD | 11/11/2015 21:39 | |

FORM I
DIESEL RANGE ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Canton | Job No.: 240-57597-1

SDG No.:

Client Sample ID: | Lab Sample ID: MB 240-205788/22-A

Matrix: Solid | Lab File ID: P5B11008.D

Analysis Method: 8015C | Date Collected:

Extraction Method: 3540C | Date Extracted: 11/09/2015 09:40

Sample wt/vol: 30(g) | Date Analyzed: 11/10/2015 16:59

Con. Extract Vol.: 5(mL) | Dilution Factor: 1

Injection Volume: 1(uL) | GC Column: ID:

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 206105 | Units: mg/Kg

| CAS NO. | COMPOUND NAME | RESULT | Q | RL | MDL |
|---|---|---|---|---|---|
| STL00141 | Total Petroleum Hydrocarbons (C10-C32) | 7.82 | J | 17 | 6.4 |

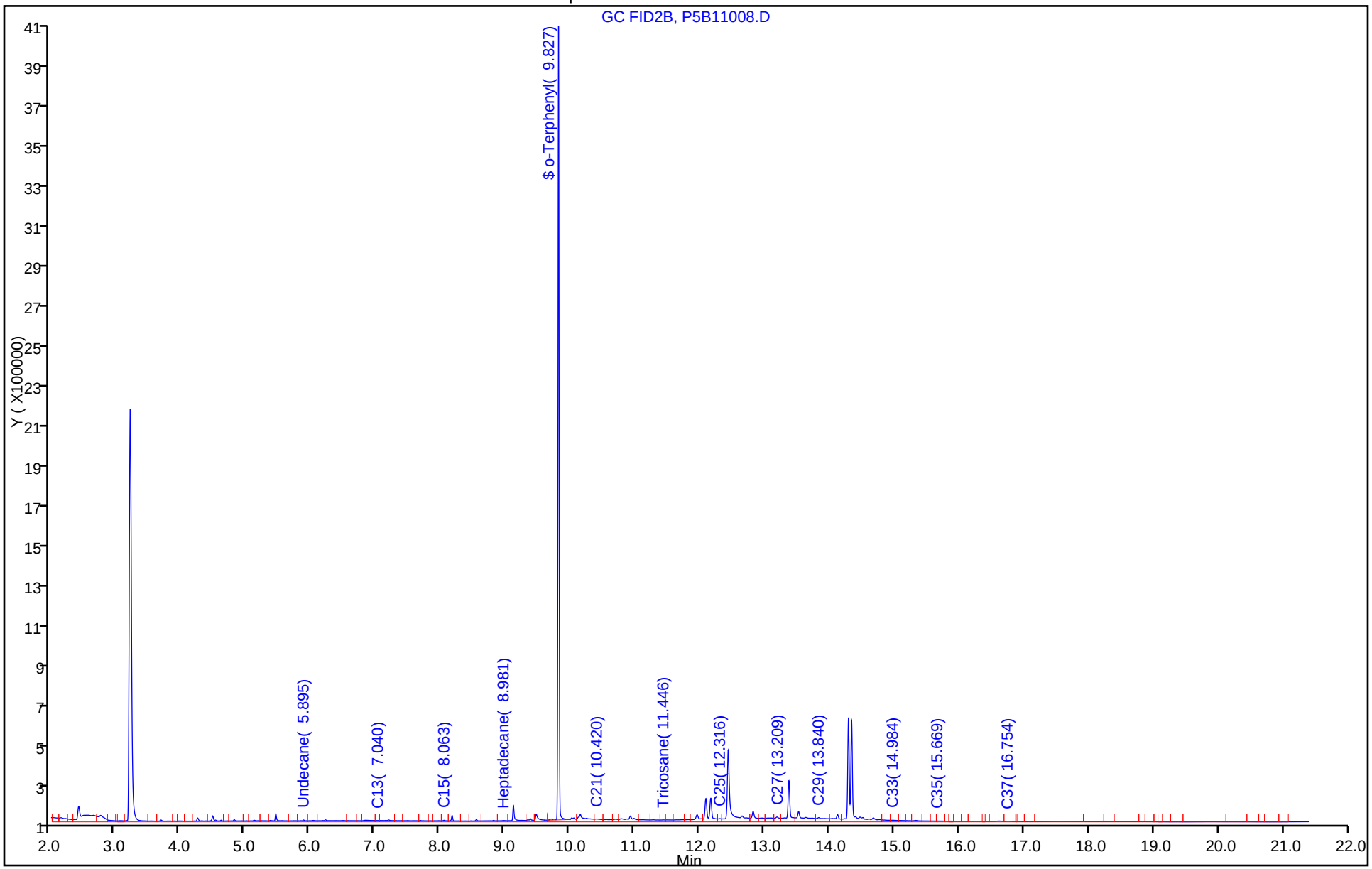
| CAS NO. | SURROGATE | %REC | Q | LIMITS |
|---|---|---|---|---|
| 84-15-1 | o-Terphenyl (Surr) | 56 | | 40-160 |

Report Date: 11-Nov-2015 11:01:18 Chrom Revision: 2.2 08-Oct-2015 07:17:48

TestAmerica Canton

Data File: \\ChromNA\Canton\ChromData\A2HP5R\20151110-49111.b\P5B11008.D

| | | | | | |
|---|---|---|---|---|---|
| Injection Date: | 10-Nov-2015 16:59:33 | Instrument ID: | A2HP5R | Operator ID: | 402582 |
| Lims ID: | MB 240-205788/22-A | | | Worklist Smp#: | 8 |
| Client ID: | | | | | |
| Injection Vol: | 1.0 ul | Dil. Factor: | 1.0000 | ALS Bottle#: | 58 |
| Method: | TPH5R | Limit Group: | GC 8015C TPH ICAL | | |



FORM III
DIESEL RANGE ORGANICS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11009.D

Lab ID: LCS 240-205788/23-A Client ID:

| COMPOUND | SPIKE ADDED (mg/Kg) | LCS CONCENTRATION (mg/Kg) | LCS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|
| Diesel | 84.0 | 53.7 | 64 | 52-120 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
DIESEL RANGE ORGANICS MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11113.D

Lab ID: 240-57597-15 MS Client ID: F-VB-UST-DUP MS

| COMPOUND | SPIKE ADDED (mg/Kg) | SAMPLE CONCENTRATION (mg/Kg) | MS CONCENTRATION (mg/Kg) | MS % REC | QC LIMITS REC | # |
|---|---|---|---|---|---|---|
| Diesel | 99.5 | 7.5 U | 75.1 | 75 | 10-124 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM III
DIESEL RANGE ORGANICS MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Matrix: Solid Level: Low Lab File ID: P5B11114.D

Lab ID: 240-57597-15 MSD Client ID: F-VB-UST-DUP MSD

| COMPOUND | SPIKE ADDED (mg/Kg) | MSD CONCENTRATION (mg/Kg) | MSD % REC | % RPD | QC LIMITS RPD | QC LIMITS REC | # |
|---|---|---|---|---|---|---|---|
| Diesel | 97.6 | 77.9 | 80 | 4 | 40 | 10-124 | |

# Column to be used to flag recovery and RPD values

FORM III 8015C

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: STD3 240-205608/10  Date Analyzed: 11/06/2015 17:59

Instrument ID: A2HP5R  GC Column:  ID: ()

Lab File ID (Standard): P5B10610.D  Heated Purge: (Y/N) N

Calibration ID: 31386

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| INITIAL CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| STD3 240-205608/10 ICRT | | 11/06/2015 17:59 | P5B10610.D | 9.83 | | |

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206105/7  Date Analyzed: 11/10/2015 16:28

Instrument ID: A2HP5R  GC Column:  ID: ()

Lab File ID (Standard): P5B11007.D  Heated Purge: (Y/N) N

Calibration ID: 31386

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCV 240-206105/7 CCVRT | | 11/10/2015 16:28 | P5B11007.D | 9.83 | | |
| MB 240-205788/22-A | | 11/10/2015 16:59 | P5B11008.D | 9.83 | | |
| LCS 240-205788/23-A | | 11/10/2015 17:31 | P5B11009.D | 9.83 | | |
| 240-57597-1 | F-VB-UST-1-14 | 11/10/2015 20:07 | P5B11014.D | 9.83 | | |
| 240-57597-2 | F-VB-UST-2-14 | 11/10/2015 20:38 | P5B11015.D | 9.83 | | |
| 240-57597-3 | F-VB-UST-3-14 | 11/10/2015 21:10 | P5B11016.D | 9.83 | | |
| 240-57597-4 | F-VB-UST-4-14 | 11/10/2015 21:41 | P5B11017.D | 9.83 | | |
| CCV 240-206105/18 | | 11/10/2015 22:12 | P5B11018.D | 9.83 | | |
| 240-57597-5 | F-VB-UST-5-14 | 11/10/2015 22:43 | P5B11019.D | 9.83 | | |
| 240-57597-6 | F-VB-UST-6-14 | 11/10/2015 23:14 | P5B11020.D | 9.83 | | |
| 240-57597-7 | F-VB-UST-7-14 | 11/10/2015 23:46 | P5B11021.D | 9.83 | | |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 00:17 | P5B11022.D | 9.83 | | |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 00:48 | P5B11023.D | 9.83 | | |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 01:19 | P5B11024.D | 9.83 | | |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 01:50 | P5B11025.D | 9.83 | | |
| CCV 240-206105/30 | | 11/11/2015 04:56 | P5B11031.D | 9.83 | | |

OTPH = o-Terphenyl

OTPH RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

FORM VIII
DIESEL RANGE ORGANICS ANALYTICAL SEQUENCE

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Sample No.: CCV 240-206356/7  Date Analyzed: 11/11/2015 18:00

Instrument ID: A2HP5R  GC Column:  ID: ()

Lab File ID (Standard): P5B11107.D  Heated Purge: (Y/N) N

Calibration ID: 31386

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

| | | | | OTPH | | |
|---|---|---|---|---|---|---|
| | | | | RT # | | |
| CONTINUING CALIBRATION SURROGATE | | | | 9.83 | | |
| UPPER LIMIT | | | | 9.93 | | |
| LOWER LIMIT | | | | 9.73 | | |
| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | LAB FILE ID | | | |
| CCV 240-206356/7 CCVRT | | 11/11/2015 18:00 | P5B11107.D | 9.83 | | |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 19:03 | P5B11109.D | 9.83 | | |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 19:34 | P5B11110.D | 9.83 | | |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 20:06 | P5B11111.D | 9.83 | | |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 20:37 | P5B11112.D | 9.83 | | |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 21:08 | P5B11113.D | 9.83 | | |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 21:39 | P5B11114.D | 9.83 | | |
| CCV 240-206356/15 | | 11/11/2015 22:11 | P5B11115.D | 9.83 | | |

OTPH = o-Terphenyl

OTPH RT Limit = ± 0.1 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8015C

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP5R Start Date: 11/06/2015 15:50

Analysis Batch Number: 205608 End Date: 11/06/2015 19:33

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| STD1 240-205608/8 IC | | 11/06/2015 15:50 | 1 | P5B10608.D | |
| STD2 240-205608/9 IC | | 11/06/2015 17:28 | 1 | P5B10609.D | |
| STD3 240-205608/10 ICRT | | 11/06/2015 17:59 | 1 | P5B10610.D | |
| STD4 240-205608/11 IC | | 11/06/2015 18:30 | 1 | P5B10611.D | |
| STD5 240-205608/12 IC | | 11/06/2015 19:01 | 1 | P5B10612.D | |
| ICV 240-205608/13 | | 11/06/2015 19:33 | 1 | P5B10613.D | |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP5R Start Date: 11/10/2015 16:28

Analysis Batch Number: 206105 End Date: 11/11/2015 04:56

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-206105/7 CCVRT | | 11/10/2015 16:28 | 1 | P5B11007.D | |
| MB 240-205788/22-A | | 11/10/2015 16:59 | 1 | P5B11008.D | |
| LCS 240-205788/23-A | | 11/10/2015 17:31 | 1 | P5B11009.D | |
| ZZZZZ | | 11/10/2015 19:04 | 1 | | |
| 240-57597-1 | F-VB-UST-1-14 | 11/10/2015 20:07 | 1 | P5B11014.D | |
| 240-57597-2 | F-VB-UST-2-14 | 11/10/2015 20:38 | 1 | P5B11015.D | |
| 240-57597-3 | F-VB-UST-3-14 | 11/10/2015 21:10 | 1 | P5B11016.D | |
| 240-57597-4 | F-VB-UST-4-14 | 11/10/2015 21:41 | 1 | P5B11017.D | |
| CCV 240-206105/18 | | 11/10/2015 22:12 | 1 | P5B11018.D | |
| 240-57597-5 | F-VB-UST-5-14 | 11/10/2015 22:43 | 1 | P5B11019.D | |
| 240-57597-6 | F-VB-UST-6-14 | 11/10/2015 23:14 | 1 | P5B11020.D | |
| 240-57597-7 | F-VB-UST-7-14 | 11/10/2015 23:46 | 1 | P5B11021.D | |
| 240-57597-8 | F-VB-UST-8-14 | 11/11/2015 00:17 | 1 | P5B11022.D | |
| 240-57597-9 | F-VB-UST-9-14 | 11/11/2015 00:48 | 1 | P5B11023.D | |
| 240-57597-10 | F-VB-UST-10-14 | 11/11/2015 01:19 | 1 | P5B11024.D | |
| 240-57597-11 | F-VB-UST-11-14 | 11/11/2015 01:50 | 1 | P5B11025.D | |
| CCV 240-206105/30 | | 11/11/2015 04:56 | 1 | P5B11031.D | |

DIESEL RANGE ORGANICS ANALYSIS RUN LOG

Lab Name: TestAmerica Canton Job No.: 240-57597-1

SDG No.:

Instrument ID: A2HP5R Start Date: 11/11/2015 18:00

Analysis Batch Number: 206356 End Date: 11/11/2015 22:11

| LAB SAMPLE ID | CLIENT SAMPLE ID | DATE ANALYZED | DILUTION FACTOR | LAB FILE ID | COLUMN ID |
|---|---|---|---|---|---|
| CCV 240-206356/7 CCVRT | | 11/11/2015 18:00 | 1 | P5B11107.D | |
| 240-57597-12 | F-VB-UST-12-14 | 11/11/2015 19:03 | 1 | P5B11109.D | |
| 240-57597-13 | F-VB-UST-13-14 | 11/11/2015 19:34 | 1 | P5B11110.D | |
| 240-57597-14 | F-VB-UST-14-14 | 11/11/2015 20:06 | 1 | P5B11111.D | |
| 240-57597-15 | F-VB-UST-DUP | 11/11/2015 20:37 | 1 | P5B11112.D | |
| 240-57597-15 MS | F-VB-UST-DUP MS | 11/11/2015 21:08 | 1 | P5B11113.D | |
| 240-57597-15 MSD | F-VB-UST-DUP MSD | 11/11/2015 21:39 | 1 | P5B11114.D | |
| CCV 240-206356/15 | | 11/11/2015 22:11 | 1 | P5B11115.D | |

FORM VI
DIESEL RANGE ORGANICS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Canton Job No.: 240-57597-1 Analy Batch No.: 205608

SDG No.:

Instrument ID: A2HP5R GC Column: ID: Heated Purge: (Y/N) N

Calibration Start Date: 11/06/2015 15:50 Calibration End Date: 11/06/2015 19:01 Calibration ID: 31386

Calibration Files:

| LEVEL: | LAB SAMPLE ID: | LAB FILE ID: |
|---|---|---|
| Level 1 | STD1 240-205608/8 | P5B10608.D |
| Level 2 | STD2 240-205608/9 | P5B10609.D |
| Level 3 | STD3 240-205608/10 | P5B10610.D |
| Level 4 | STD4 240-205608/11 | P5B10611.D |
| Level 5 | STD5 240-205608/12 | P5B10612.D |

| ANALYTE | CF | | | | CURVE TYPE | COEFFICIENT | | | # | MIN CF | %RSD | # | MAX %RSD | R^2 OR COD | # | MIN R^2 OR COD |
|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|---|
| | LVL 1 LVL 5 | LVL 2 | LVL 3 | LVL 4 | | B | M1 | M2 | | | | | | | | |
| Diesel Range Organics [C10 - C28] | 156067 174183 | 161462 | 135888 | 139742 | Ave | | 153468.493 | | | | 10.3 | | 20.0 | | | |
| o-Terphenyl (Surr) | 213204 240558 | 209043 | 182141 | 189552 | Ave | | 206899.578 | | | | 11.1 | | 20.0 | | | |

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: ICV 240-205608/13  Calibration Date: 11/06/2015 19:33

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B10613.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 164201 | | 537 | 502 | 7.0 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206105/7  Calibration Date: 11/10/2015 16:28

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11007.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 149333 | | 487 | 500 | -2.7 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 178269 | | 34.5 | 40.0 | -13.8 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206105/18  Calibration Date: 11/10/2015 22:12

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11018.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 166560 | | 543 | 500 | 8.5 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 216799 | | 41.9 | 40.0 | 4.8 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206105/30  Calibration Date: 11/11/2015 04:56

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11031.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 176144 | | 574 | 500 | 14.8 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 226941 | | 43.9 | 40.0 | 9.7 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206356/7  Calibration Date: 11/11/2015 18:00

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11107.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 156558 | | 510 | 500 | 2.0 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 215548 | | 41.7 | 40.0 | 4.2 | 20.0 |

FORM VII
DIESEL RANGE ORGANICS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Lab Sample ID: CCV 240-206356/15  Calibration Date: 11/11/2015 22:11

Instrument ID: A2HP5R  Calib Start Date: 11/06/2015 15:50

GC Column:  ID:  Calib End Date: 11/06/2015 19:01

Lab File ID: P5B11115.D  Conc. Units: ng/uL

| ANALYTE | CURVE TYPE | AVE CF | CF | MIN CF | CALC AMOUNT | SPIKE AMOUNT | %D | MAX %D |
|---|---|---|---|---|---|---|---|---|
| Diesel Range Organics [C10 - C28] | Ave | 153468 | 166194 | | 542 | 500 | 8.3 | 20.0 |
| o-Terphenyl (Surr) | Ave | 206900 | 218951 | | 42.3 | 40.0 | 5.8 | 20.0 |

# DIESEL RANGE ORGANICS BATCH WORKSHEET

Lab Name: TestAmerica Canton  Job No.: 240-57597-1

SDG No.:

Batch Number: 205788  Batch Start Date: 11/09/15 10:35  Batch Analyst: Newsome, Yvonne D

Batch Method: 3540C  Batch End Date: 11/10/15 02:35

| Lab Sample ID | Client Sample ID | Method Chain | Basis | InitialAmount | FinalAmount | exDIESELspkw 00068 | exOTPSURRW 00006 | | |
|---|---|---|---|---|---|---|---|---|---|
| 240-57597-D-1 | F-VB-UST-1-14 | 3540C, 8015C | T | 30.49 g | 5 mL | | 1 mL | | |
| 240-57597-E-2 | F-VB-UST-2-14 | 3540C, 8015C | T | 29.69 g | 5 mL | | 1 mL | | |
| 240-57597-D-3 | F-VB-UST-3-14 | 3540C, 8015C | T | 30.32 g | 5 mL | | 1 mL | | |
| 240-57597-D-4 | F-VB-UST-4-14 | 3540C, 8015C | T | 30.31 g | 5 mL | | 1 mL | | |
| 240-57597-D-5 | F-VB-UST-5-14 | 3540C, 8015C | T | 30.01 g | 5 mL | | 1 mL | | |
| 240-57597-D-6 | F-VB-UST-6-14 | 3540C, 8015C | T | 29.79 g | 5 mL | | 1 mL | | |
| 240-57597-D-7 | F-VB-UST-7-14 | 3540C, 8015C | T | 29.99 g | 5 mL | | 1 mL | | |
| 240-57597-E-8 | F-VB-UST-8-14 | 3540C, 8015C | T | 30.45 g | 5 mL | | 1 mL | | |
| 240-57597-D-9 | F-VB-UST-9-14 | 3540C, 8015C | T | 30.38 g | 5 mL | | 1 mL | | |
| 240-57597-D-10 | F-VB-UST-10-14 | 3540C, 8015C | T | 30.04 g | 5 mL | | 1 mL | | |
| 240-57597-E-11 | F-VB-UST-11-14 | 3540C, 8015C | T | 29.60 g | 5 mL | | 1 mL | | |
| 240-57597-E-12 | F-VB-UST-12-14 | 3540C, 8015C | T | 29.52 g | 5 mL | | 1 mL | | |
| 240-57597-E-13 | F-VB-UST-13-14 | 3540C, 8015C | T | 29.73 g | 5 mL | | 1 mL | | |
| 240-57597-D-14 | F-VB-UST-14-14 | 3540C, 8015C | T | 30.43 g | 5 mL | | 1 mL | | |
| 240-57597-D-15 | F-VB-UST-DUP | 3540C, 8015C | T | 29.98 g | 5 mL | | 1 mL | | |
| 240-57597-D-15 MS | F-VB-UST-DUP | 3540C, 8015C | T | 29.63 g | 5 mL | 1 mL | 1 mL | | |
| 240-57597-D-15 MSD | F-VB-UST-DUP | 3540C, 8015C | T | 30.20 g | 5 mL | 1 mL | 1 mL | | |
| MB 240-205788/22 | | 3540C, 8015C | | 30 g | 5 mL | | 1 mL | | |
| LCS 240-205788/23 | | 3540C, 8015C | | 30 g | 5 mL | 1 mL | 1 mL | | |

| Batch Notes | |
|---|---|
| Balance ID | 1339907 |
| Na2SO4 Lot Number | 2204509 |
| Prep Solvent Lot # | 2258960 |

| Basis | Basis Description |
|---|---|
| T | Total/NA |

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

# APPENDIX F—RECEIPT AND WASTE MANIFESTS

Lockheed
Block F UST's



2684189565

CAMBRIDGE IRON & METAL CO, INC
901 SOUTH KRESSON STREET
BALTIMORE, MD 21224-4504

EL0155
$$$ELITE ENVIRONMENTAL
2323 EASTERN BLVD

MIDDLE RIVER MD 21220

Rcv:590652
Date:11/12/15
Time:15:30
Ctl:590652
Ref:

| Commodity | Description | | | | |
|---|---|---|---|---|---|
| Gross | Tare | Net | Price/UM | | Amount |
| C02WF | STEEL | | | | |
| 45280* | 36700* | 8580 | 20.00 | GT | 76.61 |
| ENV | ENVIRONMENTAL FEE | | | | |
| 1* | 0 | 1 | -10.00 | EA | -10.00 |
| Totals | | 8581 | | | 66.61 |

* INDICATES MANUAL WEIGHT ENTERED

Signature:_____

ID: MATT 360 CSH 590652 MG
Veh: 3600

**Manifest Tracking Sheet**
**Soil Remedial Action - Block F UST Area**
**Lockheed Martin Corproation Middle River Complex**
**Middle River, Maryland**


| Manifest Number | Area Tracking Number | Block | Area | Truck Number | Driver | Net Weight (tons) | Net Weight (pounds) | Date |
|---|---|---|---|---|---|---|---|---|
| **Block F UST Area** | | | | | | | | |
| 343314 | TT-BF-UST-1 | F | UST | I-8 | W. Burnette | 27.53 | 55,060 | 10/23/2015 |
| 343315 | TT-BF-UST-2 | F | UST | I-6 | J. Olney | 27.44 | 54,880 | 10/23/2015 |
| 343316 | TT-BF-UST-3 | F | UST | I-15 | B. Ashford | 27.34 | 54,680 | 10/23/2015 |
| 343317 | TT-BF-UST-4 | F | UST | I-18 | C. Cofield | 30.55 | 61,100 | 10/23/2015 |
| 343318 | TT-BF-UST-5 | F | UST | I-5 | R. Crooks | 28.45 | 56,900 | 10/23/2015 |
| 343319 | TT-BF-UST-6 | F | UST | I-3 | A. Olney | 27.27 | 54,540 | 10/23/2015 |
| 343313 | TT-BF-UST-7 | F | UST | I-3 | A. Olney | 28.02 | 56,040 | 10/23/2015 |
| 343312 | TT-BF-UST-8 | F | UST | I-5 | R. Crooks | 29.63 | 59,260 | 10/23/2015 |
| 343311 | TT-BF-UST-9 | F | UST | I-18 | C. Cofield | 33.77 | 67,540 | 10/23/2015 |
| 343310 | TT-BF-UST-10 | F | UST | I-15 | B. Ashford | 33.45 | 66,900 | 10/23/2015 |
| 343309 | TT-BF-UST-11 | F | UST | I-8 | W. Burnette | 29.41 | 58,820 | 10/23/2015 |
| 343308 | TT-BF-UST-12 | F | UST | I-6 | J. Olney | 31.86 | 63,720 | 10/23/2015 |
| 343307 | TT-BF-UST-13 | F | UST | 10 | J. Thompson | 28.23 | 56,460 | 10/23/2015 |
| 343270 | TT-BF-UST-14 | F | UST | I-15 | B. Ashford | 28.12 | 56,240 | 11/2/2015 |
| 343271 | TT-BF-UST-15 | F | UST | I-13 | S. Miller | 28.62 | 57,240 | 11/2/2015 |
| 343272 | TT-BF-UST-16 | F | UST | I-3 | A. Olney | 26.78 | 53,560 | 11/2/2015 |
| 343273 | TT-BF-UST-17 | F | UST | I-6 | J. Olney | 30.44 | 60,880 | 11/2/2015 |
| 343274 | TT-BF-UST-18 | F | UST | I-18 | C. Cofield | 26.78 | 53,560 | 11/2/2015 |
| 343275 | TT-BF-UST-19 | F | UST | I-10 | D. Whalen | 31.41 | 62,820 | 11/2/2015 |
| 343276 | TT-BF-UST-20 | F | UST | I-15 | B. Ashford | 28.81 | 57,620 | 11/2/2015 |
| 343277 | TT-BF-UST-21 | F | UST | I-3 | A. Olney | 26.18 | 52,360 | 11/2/2015 |
| 343278 | TT-BF-UST-22 | F | UST | I-13 | S.Miller | 27.58 | 55,160 | 11/2/2015 |
| 343279 | TT-BF-UST-23 | F | UST | I-18 | C. Cofield | 27.23 | 54,460 | 11/2/2015 |
| 343280 | TT-BF-UST-24 | F | UST | I-6 | J. Olney | 28.22 | 56,440 | 11/2/2015 |
| 343281 | TT-BF-UST-25 | F | UST | I-10 | D. Whalen | 28.10 | 56,200 | 11/2/2015 |
| 343282 | TT-BF-UST-26 | F | UST | I-10 | J. Thompson | 30.32 | 60,640 | 11/2/2015 |
| 343283 | TT-BF-UST-27 | F | UST | PG-2 | J. Covington | 28.07 | 56,140 | 11/2/2015 |
| 343284 | TT-BF-UST-28 | F | UST | I-3 | A. Olney | 26.06 | 52,120 | 11/2/2015 |

**Manifest Tracking Sheet**
**Soil Remedial Action - Block F UST Area**
**Lockheed Martin Corproation Middle River Complex**
**Middle River, Maryland**


| Manifest Number | Area Tracking Number | Block | Area | Truck Number | Driver | Net Weight (tons) | Net Weight (pounds) | Date |
|---|---|---|---|---|---|---|---|---|
| **Block F UST Area (Continued)** | | | | | | | | |
| 343285 | TT-BF-UST-29 | F | UST | I-5 | R. Crooks | 23.99 | 47,980 | 11/3/2015 |
| 343286 | TT-BF-UST-30 | F | UST | I-7 | C. Wallace | 25.14 | 50,280 | 11/3/2015 |
| 343287 | TT-BF-UST-31 | F | UST | I-8 | W. Burnette | 31.25 | 62,500 | 11/3/2015 |
| 343288 | TT-BF-UST-32 | F | UST | I-7 | C. Wallace | 27.84 | 55,680 | 11/3/2015 |
| 343289 | TT-BF-UST-33 | F | UST | I-5 | R. Crooks | 23.08 | 46,160 | 11/3/2015 |
| 343290 | TT-BF-UST-34 | F | UST | I-3 | A. Olney | 24.57 | 49,140 | 11/3/2015 |
| 343291 | TT-BF-UST-35 | F | UST | I-8 | W. Burnette | 28.26 | 56,520 | 11/3/2015 |
| 343292 | TT-BF-UST-36 | F | UST | 10 | J. Thompson | 26.18 | 52,360 | 11/3/2015 |
| 343293 | TT-BF-UST-37 | F | UST | I-18 | C. Cofield | 32.15 | 64,300 | 11/4/2015 |
| 343294 | TT-BF-UST-38 | F | UST | I-7 | C. Wallace | 26.39 | 52,780 | 11/4/2015 |
| 343295 | TT-BF-UST-39 | F | UST | 10 | J. Thompson | 26.43 | 52,860 | 11/4/2015 |
| 343296 | TT-BF-UST-40 | F | UST | I-3 | A. Olney | 25.58 | 51,160 | 11/5/2015 |
| 343297 | TT-BF-UST-41 | F | UST | I-11 | J. McGill | 25.00 | 50,000 | 11/5/2015 |
| 343298 | TT-BF-UST-42 | F | UST | I-8 | W. Burnette | 29.33 | 58,660 | 11/5/2015 |
| 343299 | TT-BF-UST-43 | F | UST | I-20 | T. Tucker | 31.14 | 62,280 | 11/5/2015 |
| 343300 | TT-BF-UST-44 | F | UST | I-22 | J. Olley | 29.60 | 59,200 | 11/5/2015 |
| 343301 | TT-BF-UST-45 | F | UST | I-15 | B. Ashford | 30.14 | 60,280 | 11/5/2015 |
| 343302 | TT-BF-UST-46 | F | UST | I-11 | J. McGill | 25.82 | 51,640 | 11/5/2015 |
| 343303 | TT-BF-UST-47 | F | UST | I-8 | W. Burnette | 30.00 | 60,000 | 11/5/2015 |
| 343304 | TT-BF-UST-48 | F | UST | I-20 | T. Tucker | 32.32 | 64,640 | 11/5/2015 |
| 343305 | TT-BF-UST-49 | F | UST | I-15 | B. Ashford | 29.50 | 59,000 | 11/5/2015 |
| 343306 | TT-BF-UST-50 | F | UST | I-22 | J. Olley | 33.16 | 66,320 | 11/5/2015 |
| 317194 | TT-BF-UST-51 | F | UST | I-9 | M. Anderson | 29.86 | 59,720 | 11/16/2015 |
| 317195 | TT-BF-UST-52 | F | UST | I-15 | B. Ashford | 29.33 | 58,660 | 11/16/2015 |
| 317196 | TT-BF-UST-53 | F | UST | I-3 | A. Olney | 32.23 | 64,460 | 11/16/2015 |
| 317197 | TT-BF-UST-54 | F | UST | I-6 | J. Olney | 33.07 | 66,140 | 11/16/2015 |
| 317198 | TT-BF-UST-55 | F | UST | I-2 | B. Banks | 30.31 | 60,620 | 11/16/2015 |
| 317201 | TT-BF-UST-56 | F | UST | I-4 | A. Bacon | 28.07 | 56,140 | 11/16/2015 |

**Manifest Tracking Sheet**
**Soil Remedial Action - Block F UST Area**
**Lockheed Martin Corproation Middle River Complex**
**Middle River, Maryland**


| Manifest Number | Area Tracking Number | Block | Area | Truck Number | Driver | Net Weight (tons) | Net Weight (pounds) | Date |
|---|---|---|---|---|---|---|---|---|
| **Block F UST Area (Continued)** | | | | | | | | |
| 317199 | TT-BF-UST-57 | F | UST | I-15 | B. Ashford | 28.99 | 57,980 | 11/16/2015 |
| 317200 | TT-BF-UST-58 | F | UST | I-9 | M. Anderson | 26.71 | 53,420 | 11/16/2015 |
| 338192 | TT-BF-UST-59 | F | UST | I-3 | A. Olney | 30.80 | 61,600 | 11/16/2015 |
| 347857 | TT-BF-UST-60 | F | UST | I-4 | A. Bacon | 26.51 | 53,020 | 11/16/2015 |
| 347858 | TT-BF-UST-61 | F | UST | I-6 | J. Olney | 29.89 | 59,780 | 11/16/2015 |
| 347860 | TT-BF-UST-62 | F | UST | I-2 | B. Banks | 28.26 | 56,520 | 11/16/2015 |
| 342584 | TT-BF-UST-63 | F | UST | I-8 | W. Burnette | 33.39 | 66,780 | 11/23/2015 |
| 342585 | TT-BF-UST-64 | F | UST | I-15 | B. Ashford | 26.62 | 53,240 | 11/23/2015 |
| 342586 | TT-BF-UST-65 | F | UST | 10 | J. Thompson | 13.85 | 27,700 | 11/23-24/2015 |
| | | | | | **Block F UST Total** | **1850.43** | **3,700,860** | |

Note:
Soil Safe Waste Approval # W5-4194

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Log Number 29.96
TT-BE-UST-51

317194

Generator Name Lockheed Martin Corp
Address 195 Chesapeake Park Plaza Rd
Middle River MD 21220
Phone No. 410-656-4012

Shipping Location Lockheed Martin, Middle River
Address 195 Chesapeake Park Plaza
Middle River 21220
Phone No.

Approval Number WS 4194

Description of Material
Non-Regulated Petroleum Contaminated Soil
Non DOT/RCRA Regulated

GROSS
TARE
NET
TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Agent for Lockheed Martin
Generator Authorized Agent Name: Travis R Jefferson — Signature — Shipment Date: 11-16-15

### TRANSPORTER

Transporter Name: I.O.J
Address:

Driver Name (Print): Michael Anderson
Vehicle License No./State: 226868
Truck Number: I-9

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature — Shipment Date: 11.16.15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature — Delivery Date: 11.16.15

### DESTINATION

Site Name: Brandywine
Phone No.: 301-782-3036
Address: 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility  Green - Facility  Yellow - Generator  Pink - Broker  Goldenrod - Contractor  Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number: TT-BF-UST-52 29.33

317195

### GENERATOR

Generator Name: Lockheed Martin Corp. Shipping Location: Lockheed Martin Middle

Address: 195 Chesapeake Park Plaza Rd Address: 195 Chesapeake Park Pl

Middle River, Md 21220 Middle River Md 21220

Phone No.: 610-626 4012 Phone No.:

Approval Number: WS-4494

Description of Material: Non-Regulated Petroleum Contaminated Soil; Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. Ferguson for Lockheed Martin — 11-16-15

Generator Authorized Agent Name — Signature — Shipment Date

### TRANSPORTER

Transporter Name: LOS Trucking Driver Name (Print): B Ashford

Address: 3540 Crain Hwy Vehicle License No./State: 246 E96

Bowie Md Truck Number: L-15

I hereby certify that the above named material was picked up at the generator site listed above.

B Ashford 11-16-15

Driver Signature — Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

B Ashford 11-16-15

Driver Signature — Delivery Date

### DESTINATION

Site Name: Brandywine Phone No.: 301-782-3036

Address: 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number: TT-BF-051-53 3278

317196

### GENERATOR

Generator Name: Lockheed Martin Corp. | Shipping Location: Lockheed Martin Corp.

Address: 195 Chesapeake Park Plaza, Middle River MD 21220 | Address: 195 Chesapeake Park Plaza, Middle River MD 21220

Phone No.: 410-456-4013 | Phone No.:

Approval Number: WS-4494

Description of Material: Non-Regulated Petroleum Contaminated Soil; Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Tracy P. Cannon, Agent for Lockheed Martin | Signature | 11-16-15

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name: T O J | Driver Name (Print): A Only

Address: Bowie MD | Vehicle License No./State: 21CD19

Truck Number: 3-3

I hereby certify that the above named material was picked up at the generator site listed above.

A Only | 11/16

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

A Only | 11/16

Driver Signature | Delivery Date

### DESTINATION

Site Name: Brandywine | Phone No.: 301-782-3036

Address: 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility | Green - Facility | Yellow - Generator | Pink - Broker | Goldenrod - Contractor | Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number: TT-BF-UST-34 33.07

317197

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name: Lockheed Martin Corp
Address: 195 Chesapeake Park Plaza Rd
Middle River MD 21220
Phone No.: 410-656-4010

Shipping Location: Lockheed Martin Corp
Address: Chesapeake Park Plaza Rd
Middle River MD 21220
Phone No.:

Approval Number: WS-4194

Description of Material:
Non-Regulated Petroleum Contaminated Soil
Non DOT/RCRA Regulated

GROSS
TARE
NET
TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James R. Ferguson Lockheed Martin — 11-16-15

Generator Authorized Agent Name — Signature — Shipment Date

### TRANSPORTER

Transporter Name: I. O.J
Address: Bowie MD
Driver Name (Print): J. Only
Vehicle License No./State: 212 rd 46
Truck Number: I.C

I hereby certify that the above named material was picked up at the generator site listed above.

J. Only — 11/16/15

Driver Signature — Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

J. Only — 11/16/15

Driver Signature — Delivery Date

### DESTINATION

Site Name: Brandywine
Phone No.: 301-782-3036
Address: 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility  Green - Facility  Yellow - Generator  Pink - Broker  Goldenrod - Contractor  Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number: TT-BF-UST-55 30.3

317198

### GENERATOR

Generator Name Lockwood Marcell Corp  Shipping Location Lockwood Marcell Corp

Address 195 Hampton Park Plaza #1  Address 195 Hampton Park Plaza #1

White Plains MD 20695  White Plains MD 20695

Phone No. 610-656-4013  Phone No.

Approval Number: WS-4194

Description of Material: Non-Regulated Petroleum Contaminated Soil; Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James D Lucas Agent for Lockwood Marcell  11-16-15

Generator Authorized Agent Name  Signature  Shipment Date

### TRANSPORTER

Transporter Name IOJ  Driver Name (Print) Bill Banks

Address Bowie, MD  Vehicle License No./State 205 DD 8[illegible]

Truck Number 2

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature  11/16/15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature  11/16/15 Delivery Date

### DESTINATION

Site Name Brandywine  Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent  Signature  Receipt Date

White - Facility  Green - Facility  Yellow - Generator  Pink - Broker  Goldenrod - Contractor  Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 11:20

TT-BF-UST-57 28.99

317199

### GENERATOR

Generator Name Lockheed Martin Corp

Address 175 Chesapeake Park Plaza Rd

Millers Run MD 21220

Phone No. 410-656-4012

Shipping Location Lockheed Martin Corp

Address 175 Chesapeake Park Plaza Rd

Millers Run MD 21220

Phone No.

Approval Number WS-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. Ferguson Agent for Lockheed Martin JPF 11-16-15

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name T OS Trucking

Address 3540 Crain Hwy

Bowie Md

Driver Name (Print) B Ashford

Vehicle License No./State 246396

Truck Number 1-15

I hereby certify that the above named material was picked up at the generator site listed above.

B Ashford 11-16-15

Driver Signature Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

B Ashford 11-16-15

Driver Signature Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 11:30

TT-BF-UST-58 26.71

317200

### GENERATOR

Generator Name [illegible] Shipping Location [illegible]

Address [illegible] Address [illegible]

[illegible]

Phone No. [illegible] Phone No. _____

Approval Number WS-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

[illegible] 11-16-15

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name T.O.J Driver Name (Print) Michael Anderson

Address _____ Vehicle License No./State 226566

_____ Truck Number T-9

I hereby certify that the above named material was picked up at the generator site listed above.

[illegible] 11.16.15

Driver Signature Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

[illegible] 11.16.15

Driver Signature Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

_____

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number: TT-BF-051-56 28.07

317201

### GENERATOR

Generator Name: Lockheed Martin Corp

Address: 195 Chesapeake Park Plaza Rd
Middle River MD 21220

Phone No.: 410-686-4012

Shipping Location: Lockheed Martin Bldg 3

Address: 195 Chesapeake Park Plaza Rd
Middle River MD 21220

Phone No.:

Approval Number: 054174

Description of Material:
Non-Regulated Petroleum
Contaminated Soil
Non DOT/RCRA Regulated

GROSS
TARE
NET
TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Thomas P. Ferguson Agent for Lockheed Martin — 11-16-15

Generator Authorized Agent Name — Signature — Shipment Date

### TRANSPORTER

Transporter Name: E&S

Address: Bowie MD

Driver Name (Print): Burns

Vehicle License No./State: 217-45

Truck Number: E4

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature — Shipment Date: 11-16-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature — Delivery Date: 11-16-15

### DESTINATION

Site Name: Brandywine — Phone No.: 301-782-3036

Address: 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility  Green - Facility  Yellow - Generator  Pink - Broker  Goldenrod - Contractor  Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 3080

TI-96-051-59

338192

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number WS-[illegible]

4/14

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name [illegible]

Address [illegible]

Driver Name (Print) [illegible]

Vehicle License No./State [illegible]

Truck Number [illegible]

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 11-16-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 11-16-15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 33.39

TT BF WJ -63

342584

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number WS-4168 94

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Agent for Lockheed Martin

Generator Authorized Agent Name | Signature | 11-23-15 Shipment Date

### TRANSPORTER

Transporter Name JJT

Address

Burtis MD

Driver Name (Print) Wendell Burtis

Vehicle License No./State 336 567

Truck Number 5-8

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | 11-23-15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | 11-23-15 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number: 2662 TT-BF-UST-64

342585

### GENERATOR

Generator Name: Lockheed Martin Corp
Address: 195 Chesapeake Park Plaza Rd
Middle River, MD 21220
Phone No.: (610) 656-4012

Shipping Location: Lockheed Martin Middle River
Address: 195 Chesapeake Park Plaza Road
Middle River, MD 21220
Phone No.:

Approval Number: W5-410

Description of Material: Non-Regulated Petroleum Contaminated Soil; Non DOT/RCRA Regulated

GROSS
TARE
NET
TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James R Ferguson Agent for Lockheed Martin — 11-23-15

Generator Authorized Agent Name — Signature — Shipment Date

### TRANSPORTER

Transporter Name: IOS Inc.
Address: 3540 Crain Hwy
Bowie Md
Driver Name (Print): B. Ashford
Vehicle License No./State: 246-596
Truck Number: F-15

I hereby certify that the above named material was picked up at the generator site listed above.

B Ashford — 11-23-15

Driver Signature — Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

B Ashford — 11-23-15

Driver Signature — Delivery Date

### DESTINATION

Site Name: Brandywine
Phone No.: 301-782-3036
Address: 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility  Green - Facility  Yellow - Generator  Pink - Broker  Goldenrod - Contractor  Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 13.85
TT-BF - UST-65

342586

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd
Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road
Middle River, MD 21220

Phone No.

Approval Number W5-41[illegible] 94

Description of Material

Non-Regulated Petroleum
Contaminated Soil

Non DOT/RCRA Regulated

GROSS
TARE
NET
TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James R. Kenyon Agent for Lockheed Martin — 11-23-15

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name Red Top

Address

Driver Name (Print) J. Thompson

Vehicle License No./State 260 240

Truck Number 10

I hereby certify that the above named material was picked up at the generator site listed above.

J. THOMPSON 11-23-15

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

J. THOMPSON 11-23-15

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility | Green - Facility | Yellow - Generator | Pink - Broker | Goldenrod - Contractor | Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 07:10

11-2-15-11

343270

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | 11-2-15 Shipment Date

### TRANSPORTER

Transporter Name [illegible] trucking

Address 3510 [illegible] Rd

[illegible] Md

Driver Name (Print) [illegible]

Vehicle License No./State 746-596

Truck Number 1-15

I hereby certify that the above named material was picked up at the generator site listed above.

[illegible] 11-2-15

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

[illegible] 11-2-15

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 07.20
FT-BF-051-15

343271

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp
Address 195 Chesapeake Park Plaza Rd
Middle River, MD 21220
Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River
Address 195 Chesapeake Park Plaza Road
Middle River, MD 21220
Phone No.

Approval Number W5-4194

Description of Material
Non-Regulated Petroleum Contaminated Soil
Non DOT/RCRA Regulated

GROSS
TARE
NET
TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P Larson Lockheed Martin — 11-2-15
Generator Authorized Agent Name — Signature — Shipment Date

### TRANSPORTER

Transporter Name I.O.J.
Address 3540 Crain Hwy # 208
Bowie, MD 20716

Driver Name (Print) S. Miller
Vehicle License No./State 245 74MD
Truck Number I-13

I hereby certify that the above named material was picked up at the generator site listed above.

[signature] 11-2-15
Driver Signature — Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

[signature] 11-2-15
Driver Signature — Delivery Date

### DESTINATION

Site Name Brandywine
Phone No. 301-782-3036
Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility  Green - Facility  Yellow - Generator  Pink - Broker  Goldenrod - Contractor  Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 07:26

FT-GE-UST-16

343272

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name J. O. J.

Address Bowie MD

Driver Name (Print) Comp Name Only

Vehicle License No./State 211219

Truck Number J-3

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine

Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility | Green - Facility | Yellow - Generator | Pink - Broker | Goldenrod - Contractor | Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 07:33

## NON-HAZARDOUS MATERIAL MANIFEST

343273

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. Lawson Lockheed Martin

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name E C J

Address

Bowie MD

Driver Name (Print) J. Cisko

Vehicle License No./State 212 CL46

Truck Number 26

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 11/3/15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 11/3/15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 07.46

11-01-151-18

343274

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P Lawson Lockheed Martin 11-2-15

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name I.O.J

Address BOWIE, MD

Driver Name (Print) C Cofield

Vehicle License No./State 253525

Truck Number I-18

I hereby certify that the above named material was picked up at the generator site listed above.

C Cofield 11.2.15

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

C Cofield 11-2-15

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 07:55

11-BF-U31 19

343275

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name IOJ

Address 3540 Crain hwy

Bowie Md

Driver Name (Print) Derrick Whalen

Vehicle License No./State 236ED89

Truck Number I-10

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 11/2/15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 11/2/15

### DESTINATION

Site Name Brandywine

Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 10:12

TT-BE-WSI-20

343276

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name Signature 11-2-15 Shipment Date

### TRANSPORTER

Transporter Name T O Trucking

Address 3610 Clay Hill

Bowie MD

Driver Name (Print) B Ballard

Vehicle License No./State 246-396

Truck Number T-15

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature 11-2-15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature 11-2-15 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 10:44

FT-BF-UST-39

343277

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Jason P. Ferguson Agent for Lockheed Martin — 11-3-15

Generator Authorized Agent Name — Signature — Shipment Date

### TRANSPORTER

Transporter Name [illegible]

Address [illegible]

Driver Name (Print) [illegible]

Vehicle License No./State [illegible]

Truck Number 1-3

I hereby certify that the above named material was picked up at the generator site listed above.

A Cruley 11/2/15

Driver Signature — Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

A Cruley 11/2/15

Driver Signature — Delivery Date

### DESTINATION

Site Name Brandywine — Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility  Green - Facility  Yellow - Generator  Pink - Broker  Goldenrod - Contractor  Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 10:54

11-BE-1151-22

343278

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Agent for

James P. Sorensen Jr. Lockheed Martin

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11.00

TT-BF-UST-23

343279

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. Farmer Lockheed Martin 11-2-15

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11:05

11-BE-141-TH

343280

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11:16

11-BF-051-35

343281

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4198

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. Ferguson for Lockheed Martin 11-3-13

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name J&S

Address 3540 Crain hwy

Bowie MD

Driver Name (Print) D Wheeler

Vehicle License No./State 236509

Truck Number 1-10

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature 11-3-13 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature 11-3-13 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number

LT-BF-1137-16

343282

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

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Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name R&D TRUCKING/IOS

Address

Driver Name (Print) JERMAINE Thompson

Vehicle License No./State 320 6.16

Truck Number I 10

I hereby certify that the above named material was picked up at the generator site listed above.

J. THOMPSON 11-2-15

Driver Signature Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

J. THOMPSON 11-2-15

Driver Signature Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 14:20

11-RI-UST-27

343283

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P Ingram Lockheed Martin 11/2/15

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name Pinnacle Group

Address Upper Marlboro

Driver Name (Print) Brad Gaines

Vehicle License No./State 24V284

Truck Number P6-2

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature 11/2/15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature 11/2/15 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 14135

343284

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 07-15

11-BF-UST-39

343285

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. [illegible] Lockheed Martin — 11-3-15

Generator Authorized Agent Name — Signature — Shipment Date

### TRANSPORTER

Transporter Name F.O.J

Address Bowie rd

Driver Name (Print) Raymond Clark

Vehicle License No./State MD2[illegible]

Truck Number F-5

I hereby certify that the above named material was picked up at the generator site listed above.

11-03-15

Driver Signature — Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

11-3-15

Driver Signature — Delivery Date

### DESTINATION

Site Name Brandywine

Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility  Green - Facility  Yellow - Generator  Pink - Broker  Goldenrod - Contractor  Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 07:25

T1-BE-USJ-Z

343286

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date 11-3-15

### TRANSPORTER

Transporter Name IUS

Address 3540 Crain Hwy

Bowie MD

Driver Name (Print) Craig Wallace

Vehicle License No./State 246D58

Truck Number I-7

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 11-3-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 11-3-15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 08:25

TT-BF-051-31

343287

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. Vinciguerra Lockheed Martin 11-3-15

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name TDJ

Address

Driver Name (Print) Wendell Puckett

Vehicle License No./State 224643

Truck Number 3-8

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature 11-3-15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature 11-3-15 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 10:49

11-BE-UST-30

343288

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W8-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 11:00

11-BF-UST-33

343289

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. Fanning for Lockheed Martin 11-3-15

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name JOJ

Address 3140 Long Meadow

Bowie Md

Driver Name (Print) R. Brooks

Vehicle License No./State 219 B 57

Truck Number 3-5

I hereby certify that the above named material was picked up at the generator site listed above.

11-3-15

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

11-3-15

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11:46

## NON-HAZARDOUS MATERIAL MANIFEST

343290

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 12,15
TT-BF-UST-36

343291

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date 11-3-15

### TRANSPORTER

Transporter Name TJJ

Address

Driver Name (Print) Wendell Burnette

Vehicle License No./State 226647

Truck Number 1-3

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 11-3-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 11-3-15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 13:00

## NON-HAZARDOUS MATERIAL MANIFEST

343292

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name [illegible] TOT

Address HAGERSTOWN MARYLAND

Driver Name (Print) JERMAINE THOMPSON

Vehicle License No./State [illegible]

Truck Number 10

I hereby certify that the above named material was picked up at the generator site listed above.

J. THOMPSON

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

J. THOMPSON

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 10150

TT-BT-US1-37

343293

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date 11-4-15

### TRANSPORTER

Transporter Name IOI

Address 3340 Liberty Highway

Boise ME

Driver Name (Print) C Canfield

Vehicle License No./State 033 EP 25

Truck Number T-19

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 11-4-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 11-4-15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 12:44

343294

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name _____ Signature _____ Shipment Date _____

### TRANSPORTER

Transporter Name _____

Address _____

_____

Driver Name (Print) _____

Vehicle License No./State _____

Truck Number _____

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature _____ Shipment Date _____

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature _____ Delivery Date _____

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent _____ Signature _____ Receipt Date _____

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 13:00

11-BF-057-27

343295

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Agent Name | Signature | 11-4-15

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name R.L. Inc

Address

Driver Name (Print) J Thompson

Vehicle License No./State 240 5 40

Truck Number 10

I hereby certify that the above named material was picked up at the generator site listed above.

J. THOMPSON 11-4-15

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

J. THOMPSON 11-4-15

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility | Green - Facility | Yellow - Generator | Pink - Broker | Goldenrod - Contractor | Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 07:20

TT-BF-USF-40

343296

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date 11.5.15

### TRANSPORTER

Transporter Name JOT

Address

Bowie Md

Driver Name (Print) N Only

Vehicle License No./State 211 5 19

Truck Number I-3

I hereby certify that the above named material was picked up at the generator site listed above.

N Only 11/5/15

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

N Only 11/5/15

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 07:10

TT-BF-UST-41

343297

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | 11-5-15 Shipment Date

### TRANSPORTER

Transporter Name JDJ

Address 3570 Old Hwy

Bowie MD

Driver Name (Print) J. M. Gill

Vehicle License No./State 240589

Truck Number J-11

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | 11-5-15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | 11-5-15 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 07:40

TI-BF-UST-42

343298

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name _____ Signature _____ Shipment Date 4-5-16

### TRANSPORTER

Transporter Name _____

Address _____

_____

Driver Name (Print) _____

Vehicle License No./State _____

Truck Number _____

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature _____ Shipment Date 4-5-16

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature _____ Delivery Date 4-6-16

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent _____ Signature _____ Receipt Date _____

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 8.15

TT-BF-VST-43

343299

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility | Green - Facility | Yellow - Generator | Pink - Broker | Goldenrod - Contractor | Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 8:30

TT - BF - UST - 44

343300

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name _____ Signature _____ 11-5-15 Shipment Date

### TRANSPORTER

Transporter Name _____

Address _____

_____

Driver Name (Print) Joe Olley

Vehicle License No./State _____

Truck Number _____

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature _____ 11-5-15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature _____ 11-5-15 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent _____ Signature _____ Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 8:45

TT-BF-UST-45

## NON-HAZARDOUS MATERIAL MANIFEST

343301

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name T&T Trucking

Address 3540 Crain Hwy

Bowie Md

Driver Name (Print) B Ashford

Vehicle License No./State 246ED96

Truck Number T-15

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 11-15-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 11-15-15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11-27

TT-BF-UST-46

343302

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name EST

Address 302 Cain Hwy

Bowie MD

Driver Name (Print) Jason McGill

Vehicle License No./State 2410 ED MD

Truck Number T-11

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 11-5-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 11-5-15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 11-27
TT-BF-UST-47

343303

### GENERATOR

Generator Name Lockheed Martin Corp
Address 195 Chesapeake Park Plaza Rd
Middle River, MD 21220
Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River
Address 195 Chesapeake Park Plaza Road
Middle River, MD 21220
Phone No.

Approval Number W5-4194

Description of Material
Non-Regulated Petroleum Contaminated Soil
Non DOT/RCRA Regulated

GROSS
TARE
NET
TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Agent for
James R Sorensen Lockheed Martin 11-5-13
Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name EQ
Address
Beaver WV
Driver Name (Print)
Vehicle License No./State 226567
Truck Number 19

I hereby certify that the above named material was picked up at the generator site listed above.

11-5-13
Driver Signature Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

11-5-13
Driver Signature Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036
Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number: TT-BF-WST-48

343304

### GENERATOR

Generator Name: Lockheed Martin Corp

Address: 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No.: (610) 656-4012

Shipping Location: Lockheed Martin Middle River

Address: 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.:

Approval Number: W5-4194

Description of Material: Non-Regulated Petroleum Contaminated Soil; Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date

### DESTINATION

Site Name: Brandywine

Phone No.: 301-782-3036

Address: 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility | Green - Facility | Yellow - Generator | Pink - Broker | Goldenrod - Contractor | Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number: TT-BF-UST-N9

343305

### GENERATOR

Generator Name: Lockheed Martin Corp
Address: 195 Chesapeake Park Plaza Rd
Middle River, MD 21220
Phone No.: (610) 656-4012

Shipping Location: Lockheed Martin Middle River
Address: 195 Chesapeake Park Plaza Road
Middle River, MD 21220
Phone No.:

Approval Number: W5-4194

Description of Material: Non-Regulated Petroleum Contaminated Soil; Non DOT/RCRA Regulated

GROSS
TARE
NET
TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name: Joes Trucking
Address: 3540 Crain Hwy
Bowie Md
Driver Name (Print): Bradford
Vehicle License No./State: 24[illegible]96
Truck Number: T-15

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature: B Gilford — Shipment Date: 11-5-15 / 11-5-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature: B Gilford — Delivery Date: 11-5-15 / 11-5-15

### DESTINATION

Site Name: Brandywine — Phone No.: 301-782-3036

Address: 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility | Green - Facility | Yellow - Generator | Pink - Broker | Goldenrod - Contractor | Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number: TT-BF-UST-50

343306

### GENERATOR

Generator Name: Lockheed Martin Corp

Address: 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No.: (610) 656-4012

Shipping Location: Lockheed Martin Middle River

Address: 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.:

Approval Number: W5-4194

Description of Material: Non-Regulated Petroleum Contaminated Soil; Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name: [illegible] Agent for Lockheed Martin

Signature

Shipment Date: 11-5-15

### TRANSPORTER

Transporter Name: TOS

Address:

Bowie Md

Driver Name (Print): T. Olky

Vehicle License No./State: 256 5 45

Truck Number: I-22

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature

Shipment Date: 11-5-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature

Delivery Date: 11-5-15

### DESTINATION

Site Name: Brandywine

Phone No.: 301-782-3036

Address: 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent

Signature

Receipt Date

White - Facility  Green - Facility  Yellow - Generator  Pink - Broker  Goldenrod - Contractor  Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number

343307

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name [illegible]

Address [illegible]

Driver Name (Print) [illegible] Thomas

Vehicle License No./State [illegible]

Truck Number 10

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 10/23/15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 10/23/15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 12.15

343308 12

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name EOT

Address

Driver Name (Print) T. Cody

Vehicle License No./State 212446

Truck Number 76

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 10/20/15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 10/20/15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11 12:30

343309

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print) Warde II Burnette

Vehicle License No./State 206 ED 67

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 10/2/15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 10/2/15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11:30

343310

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

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Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11:40

343311

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number T-18

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date 10-23-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date 10-23-15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11:19

## NON-HAZARDOUS MATERIAL MANIFEST

343312

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print) Raymond Cooks

Vehicle License No./State MD 1411 C2

Truck Number T-5

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 11:00

343313

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name

Address

Driver Name (Print)

Vehicle License No./State

Truck Number

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 9:00

343314

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4184

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P Lanahan Agent for Lockheed Martin 10/23/15

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name TDS

Address 3546 Crain Hwy

Bowie MD

Driver Name (Print) Woodell Burnette

Vehicle License No./State 22 FD61

Truck Number 18

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature 10/23/15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature 10/23/15 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 8375

11-BF-UST-2

343315

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

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James P. Ferguson for Lockheed Martin 10-23-15

Generator Authorized Agent Name Signature Shipment Date

### TRANSPORTER

Transporter Name EOT

Address Bowie MD

Driver Name (Print) J Ortiz

Vehicle License No./State 212 GA 46

Truck Number E-6

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature 10/23/15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature 10/23/15 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 8:30

343316

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name Signature Shipment Date 10-23-15

### TRANSPORTER

Transporter Name L & S Trucking

Address 3500 Conowingo Rd

Bruce MD

Driver Name (Print)

Vehicle License No./State 2046E94

Truck Number 1-15

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature Shipment Date 10-23-15

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature Delivery Date 10-23-15

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 8:00

T1-B1 - UST-4

343317

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. Ferguson Agent for Lockheed Martin — 10-23-15

Generator Authorized Agent Name | Signature | Shipment Date

### TRANSPORTER

Transporter Name I.O.J

Address Bowie, MD

Driver Name (Print) C Cofield

Vehicle License No./State 253 E 25

Truck Number I-18

I hereby certify that the above named material was picked up at the generator site listed above.

C Cofield 10-23-15

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

C Cofield 10-23-15

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 7:45

11-05-031-5

343318

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W6-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Agent for

Generator Authorized Agent Name Signature 10-23-15 Shipment Date

### TRANSPORTER

Transporter Name T & D

Address Bowie md

Driver Name (Print) Raymond Cooks

Vehicle License No./State MD 2111S7

Truck Number I-5

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature 10-23-15 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature 10-23-15 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 7:35

T1-BL-057-6

343319

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 656-4012

Shipping Location Lockheed Martin Middle River

Address 195 Chesapeake Park Plaza Road

Middle River, MD 21220

Phone No.

Approval Number W5-4194

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

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Generator Authorized Agent Name | Signature | Shipment Date 10/23/15

### TRANSPORTER

Transporter Name I.O.I.

Address Bowie MD

Driver Name (Print) Antonio Only

Vehicle License No./State 211ED19

Truck Number 13

I hereby certify that the above named material was picked up at the generator site listed above.

A Only 10/23/15

Driver Signature | Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

A Only 10/23/15

Driver Signature | Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent | Signature | Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 11:57 TT-BF-USF-60 24.51

347857

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd.
Middle River, MD 21220

Phone No. (610) 636-4012

Shipping Location Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd.
Middle River, MD 21220

Phone No. (610) 636-4012

Approval Number WS-4198

Description of Material

Non-Regulated Petroleum
Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

James P. Gagnon for Lockheed Martin — 11-16-15

Generator Authorized Agent Name — Signature — Shipment Date

### TRANSPORTER

Transporter Name JJ

Address B
Bowie md

Driver Name (Print) A Bacon

Vehicle License No./State 217 563

Truck Number E-1

I hereby certify that the above named material was picked up at the generator site listed above.

A Bacon — 11-16-15

Driver Signature — Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

A Bacon — 11-16-15

Driver Signature — Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

## NON-HAZARDOUS MATERIAL MANIFEST

Log Number 29.89
TT-BF-UST-61

347858

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd.
Middle River, MD 21220

Phone No. (610) 636-4012

Shipping Location Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd.
Middle River, MD 21220

Phone No. (610) 636-4012

Approval Number W5-4198-4

Description of Material

Non-Regulated Petroleum
Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

**I hereby certify that the above named material does not contain free liquid as defined by 40 CFR Part 260.10 or any applicable state law, is not a hazardous waste as defined by 40 CFR Part 261 or any applicable state law, has been properly described, classified and packaged, and is in proper condition for transportation according to applicable regulations.**

Generator Authorized Agent Name Agent for Lockheed Martin Signature 11-16-13 Shipment Date

### TRANSPORTER

Transporter Name EQ

Address Baltimore MD

Driver Name (Print) J Olney

Vehicle License No./State 212546

Truck Number 2-6

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature 11/16/13 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature 11/16/13 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent Signature Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# SOIL SAFE, INC.

Log Number 22.26

IT-BL-USF-62

347860

## NON-HAZARDOUS MATERIAL MANIFEST

### GENERATOR

Generator Name Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd.

Middle River, MD 21220

Phone No. (610) 636-4012

Shipping Location Lockheed Martin Corp

Address 195 Chesapeake Park Plaza Rd

Middle River, MD 21220

Phone No. (610) 636-4012

Approval Number 4 WS-4191

Description of Material

Non-Regulated Petroleum Contaminated Soil

Non DOT/RCRA Regulated

GROSS

TARE

NET

TONNAGE

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Thomas P. Furgeson Agent for Lockheed Martin

Generator Authorized Agent Name — Signature — 11-16-13 Shipment Date

### TRANSPORTER

Transporter Name JDS

Address Bowie MD

Driver Name (Print) B. Banks

Vehicle License No./State JDS D 88

Truck Number 1-2

I hereby certify that the above named material was picked up at the generator site listed above.

Driver Signature — 11-16-13 Shipment Date

I hereby certify that the above named material was delivered without incident to the destination listed below.

Driver Signature — 11-16-13 Delivery Date

### DESTINATION

Site Name Brandywine Phone No. 301-782-3036

Address 16001 Mattawoman Drive, Brandywine, Md. 20613

I hereby certify that the above named material has been accepted and to the best of my knowledge the foregoing is true and accurate.

Name of Authorized Agent — Signature — Receipt Date

White - Facility Green - Facility Yellow - Generator Pink - Broker Goldenrod - Contractor Blue - Trucking Co.

# APPENDIX G—30-DAY NOTIFICATION FORM, REGISTRATION FORMS, AND MDE TANK REMOVAL REPORT

# MARYLAND DEPARTMENT OF THE ENVIRONMENT

Land Management Administration • Oil Control Program
1800 Washington Boulevard • Suite 620 • Baltimore Maryland 21230-1719
410-537-3442 • 800-633-6101 x3442 • 410-537-3092 (fax) • www.mde.maryland.gov

## Underground Storage System Removal/Abandonment

## 30-Day Written Notification

**Case No:** _____

☐ Facility No:_____
(check box if facility was not previously registered)

This form shall be used to notify the Department at least 30 days before beginning underground storage tank removal and/or abandonment-in-place. When fully completed, this form may be accepted as an amendment to the Notification for Underground Storage Tanks currently on file with the Department, for the removals and/or abandonments listed. New tank installations must be reported on the five-page notification (Form Number MDE/WAS/PER.012). The Department reserves the right to require Form Number MDE/WAS.PER.012, if determined necessary to properly update Department records.

(1) ***Type of facility***: (check one) ___Government X Commercial ___Farm/Nursery ___Residential (non-rental) ___Other (please specify)_____

(2) ***Type of work being performed***: (check all that apply) X Removal ___Abandonment in Place ___Temporary Closure ___Installation ___Upgrade of Existing Tank/Piping

(3) ***Date work is to be performed***: July 28, 2015 (4) ***Estimated time that work will be ready for inspection***: 10:00 am

(5) ***Insurance Information***: (check one) ___Self Insurance ___Insurance Pool ___Risk Retention Group ___Guarantee ___Letter of Credit ___Surety Bond
___Commercial Insurance: Policy No.:_____ Insurer:_____ Agent/Broker:_____ Phone:_____
___ Other Method allowed: (specify)_____

| (6) ***Contractor Information:*** | (7) ***Facility Information:*** | (8) ***Owner Information:*** |
|---|---|---|
| Elite Environmental & Petroleum Services, Inc. | Lockheed Martin Corporation Middle River Complex Tax Block F | Lockheed Martin Corporation |
| Company Name | Facility Name | Owner Name |
| P.O. Box 615 | 195 Chesapeake Park Plaza | 6801 Rockledge Drive |
| Mailing Address | Street Address | Mailing Address |
| Cape Canaveral, FL 32920 | Middle River, MD 21220 | Bethesda, MD 20817 |
| City/State/Zip | City/State/Zip | City/State/Zip |
| Lawrence Macher | Eastern Blvd. | Tom Blackman |
| Name of Contact Person | Nearest Cross Street | Contact Person at owner location (**not contractor**) |
| 443.324.1100 321.613.5267 | Jim Ferguson | 301.548.2209 |
| Telephone No. Fax No. | Name of Contact Person at Site | Telephone No. Fax No. |
| Lawrence Macher | 412.496.9283 | Michael Martin (Tetra Tech Program Manager) |
| Name of Person certified to do work | Telephone No. of Contact Person | Name/Title of person authorized to represent owner |
| MDIC- 2015-1907 exp. date 05/01/2017 | | |

## 30-DAY WRITTEN NOTIFICATION

MDE Oil Control Program

(9) ***Underground Storage Tank Information:*** ***Facility No.:*** _____

| Tank Number | Tank Capacity | Type of Product | Material of Construction Tank | Material of Construction Piping | Date Tank Last Used | Date Tank Last Tested | Pass or Failed? | Type of Test |
|---|---|---|---|---|---|---|---|---|
| 1 | 1,000 gal | unknown | unknown | unknown | 1954 | unknown | — | — |
| 2 | 1,000 gal | unknown | unknown | unknown | 1954 | unknown | — | — |
| 3 | 1,000 gal | unknown | unknown | unknown | 1954 | unknown | — | — |
| 4 | 1,000 gal | unknown | unknown | unknown | 1954 | unknown | — | — |
| 5 | 1,000 gal | unknown | unknown | unknown | 1954 | unknown | — | — |
| 6 | 1,000 gal | unknown | unknown | unknown | 1954 | unknown | — | — |
| 7 | 1,000 gal | unknown | unknown | unknown | 1954 | unknown | — | — |
| | | | | | | | | |
| | | | | | | | | |

(10) ***Are there additional underground storage tanks at this facility not listed above?*** ____Yes __X__No

(11) ***Certification***:

I certify, under penalty of law, that I have personally examined and am familiar with the information submitted in this and all attached documents. Based on my inquiry of those individuals immediately responsible for obtaining the information, I believe that the information submitted is true, accurate and complete. I understand this form may not be accepted by the Oil Control Program if the information is incomplete. (Complete items 1 through 11)

***Signature of UST Owner/ Authorized Owner Representative:*** _____ ***Title:*** Tetra Tech Program Manager ***Date:*** 5/29/15
(as listed in section 8 of this form)

**Notice: Collection of Personal Records – State Government Article § 10-624**

This Notice is provided pursuant to § 10-624 of the State Government Article of the Maryland Code. The personal information requested on this form is intended to be used in processing your application. Failure to provide the information requested may result in your application not being processed. You have the right to inspect, amend, or correct this form. The Maryland Department of the Environment ("MDE") is a public agency and subject to the Maryland Public Information Act. This form may be made available on the Internet via MDE's website and is subject to inspection or copying, in whole or in part, by the public and other governmental agencies, if not protected by federal or State law.

# MARYLAND DEPARTMENT OF THE ENVIRONMENT

1800 Washington Boulevard, Suite 620 • Baltimore Maryland 21230-1719
(410) 537-3442 • 1-800-633-6101 • http:// www.mde.maryland.gov

P1

LAND MANAGEMENT ADMINISTRATION
Oil Control Program

Case #: 16-0084BA
☐INITIAL / ☒FOLLOW-UP

## Tank Removal / Abandonment

Site Name: BLOCK F - LOCHHEED MARTIN CORP
Address: 195 CHESAPEAK PARK PLAZA MIDDLE RIVER, MD 21220

Date: 11-4-2015
Facility ID #: 13189

1. (a) 7 Number of UST's removed (b) 0 Number of UST's abandoned-in-place (c) 1 Number of UST's remaining on-site.
2. Has an environmental assessment been completed? ☐YES ☒NO

| Tank | Product | Age (years) | Size (gallons) | Tank Construction | Perforations | | Piping Construction | Disposal Site |
|---|---|---|---|---|---|---|---|---|
| | | | | | Tank | Piping | | |
| 1 | Aviation | >25 | 1000 | Bare Steel | Yes ☐ No ☒ | Yes ☐ No ☐ | Bare Steel | CAMBRIDGE |
| 2 | Aviation | >25 | 1000 | Bare Steel | Yes ☐ No ☒ | Yes ☐ No ☐ | | Cambridge |
| 3 | Aviation | >25 | 1000 | Bare Steel | Yes ☐ No ☒ | Yes ☐ No ☐ | | Cambridge |
| 4 | Aviation | >25 | 1000 | Bare Steel | Yes ☐ No ☒ | Yes ☐ No ☐ | | Cambridge |
| 5 | Aviation | >25 | 1000 | Bare Steel | Yes ☐ No ☒ | Yes ☐ No ☐ | | Cambridge |
| 6 | Aviation | >25 | 1000 | Bare Steel | Yes ☐ No ☒ | Yes ☐ No ☐ | | Cambridge |

3. Has piping been properly abandoned? ☐YES ☒NO ☐UNKNOWN
4. Has vent risers been removed? ☐YES ☐NO N/A
5. Has all liquid been removed from UST(s)? ☒YES ☐NO
6. Is explosion meter on site? ☒YES ☐NO
7. Have UST(s) been purged of explosive or combustible vapors? ☒YES ☐NO
   (must confirm less than 10% LEL with explosion meter)
8. Is groundwater contaminated? ☐YES ☐NO ☒NOT DETECTABLE AT THIS TIME
9. Is soil contaminated? (if yes, type of product:_____) ☐YES ☐NO ☒NOT DETECTABLE AT THIS TIME
10. Were contaminated soils removed? ☒YES ☐NO
    (If YES, complete *Contaminated Soil Removal Form*; If NO, describe in item 18)
11. Was soil field screened? ☒YES ☐NO
    Tank: max. units 4 at overburden Piping: max. units _____ at _____
12. Are domestic well(s) on site? ☐YES ☒NO
    Is sampling required? (If YES, list EPA method in item 14) ☐YES ☐NO

13. **ACTIONS REQUIRED, IMMEDIATELY, OF THE OWNER BY THIS ADMINISTRATION:**
☐STOP OPERATIONS ☐PUMP OUT LIQUIDS ☐CONTAIN AND CLEAN UP SPILL
☐OTHER: _____

14. **ACTIONS REQUIRED, WITHIN 45 DAYS, OF THE OWNER BY THIS ADMINISTRATION:**
☒SUBMIT ALL TANK REMOVAL/ABANDONMENT DOCUMENTATION INCLUDING: UST CLOSURE REPORT as per guidance document.
☒PROPERLY ABANDON PIPING IN COMPLIANCE WITH COMAR 26.10.10.02 B.(2) (removed unless otherwise directed)
☒REMOVE VENT PIPE RISER (S)
☐_____ MONITORING WELL (S) REQUIRED IN PETROLEUM IMPACTED AREA (S) DESCRIBED IN ITEM 18
☐COMPLETE AN ENVIRONMENTAL ASSESSMENT IN COMPLIANCE WITH COMAR 26.10._____(submit two copies)
☒SUBMIT SOIL ANALYTICAL RESULTS:
**EPA METHOD:** ☒8015B GRO/DRO ☒8270 (SVOCs) ☒8260 (VOCs)
☒OTHER:PCBS EPA 8081/8082
☐SUBMIT GROUNDWATER ANALYTICAL RESULTS:
**EPA METHOD** ☐8015B GRO/DRO ☐8270 (SVOCs) ☐8260 (VOCs) ☐524.2 (VOCs)
☐OTHER _____
☒SUBMIT SOIL DISPOSAL RECEIPT
☒SUBMIT TANK DISPOSAL RECEIPT

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# MARYLAND DEPARTMENT OF THE ENVIRONMENT

1800 Washington Boulevard, Suite 620 • Baltimore Maryland 21230-1719
(410) 537-3442 • 1-800-633-6101 • http:// www.mde.maryland.gov

LAND MANAGEMENT ADMINISTRATION
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1. (a) 7 Number of UST's removed (b) 0 Number of UST's abandoned-in-place (c) 1 Number of UST's remaining on-site.
2. Has an environmental assessment been completed? ☐YES ☒NO

| Tank | Product | Age (years) | Size (gallons) | Tank Construction | Perforations | | Piping Construction | Disposal Site |
|---|---|---|---|---|---|---|---|---|
| | | | | | Tank | Piping | | |
| 7 | Aviation | >25 | 1000 | Bare Steel | Yes ☐ No ☒ | Yes ☐ No ☐ | Bare Steel | CAMBRIDGE |
| | | | | | Yes ☐ No ☒ | Yes ☐ No ☐ | | |
| | | | | | Yes ☐ No ☒ | Yes ☐ No ☐ | | |
| | | | | | Yes ☐ No ☒ | Yes ☐ No ☐ | | |
| | | | | | Yes ☐ No ☒ | Yes ☐ No ☐ | | |
| | | | | | Yes ☐ No ☒ | Yes ☐ No ☐ | | |

3. Has piping been properly abandoned? ☐YES ☒NO ☐UNKNOWN
4. Has vent risers been removed? ☐YES ☐NO NA
5. Has all liquid been removed from UST(s)? ☒YES ☐NO
6. Is explosion meter on site? ☒YES ☐NO
7. Have UST(s) been purged of explosive or combustible vapors? ☒YES ☐NO
   (must confirm less than 10% LEL with explosion meter)
8. Is groundwater contaminated? ☐YES ☐NO ☒NOT DETECTABLE AT THIS TIME
9. Is soil contaminated? (if yes, type of product:_____) ☐YES ☐NO ☒NOT DETECTABLE AT THIS TIME
10. Were contaminated soils removed? ☒YES ☐NO
    (If YES, complete *Contaminated Soil Removal Form*; If NO, describe in item 18)
11. Was soil field screened? ☒YES ☐NO
    Tank: max. units 4 at overburden Piping: max. units _____ at _____
12. Are domestic well(s) on site? ☐YES ☒NO
    Is sampling required? (If YES, list EPA method in item 14) ☐YES ☐NO
13. **ACTIONS REQUIRED, IMMEDIATELY, OF THE OWNER BY THIS ADMINISTRATION**:
    ☐STOP OPERATIONS ☐PUMP OUT LIQUIDS ☐CONTAIN AND CLEAN UP SPILL
    ☐OTHER: _____

14. **ACTIONS REQUIRED, WITHIN 45 DAYS, OF THE OWNER BY THIS ADMINISTRATION**:
    ☒SUBMIT ALL TANK REMOVAL/ABANDONMENT DOCUMENTATION INCLUDING: UST CLOSURE REPORT as per guidance document.
    ☒PROPERLY ABANDON PIPING IN COMPLIANCE WITH COMAR 26.10.10.02 B.(2) (removed unless otherwise directed)
    ☐REMOVE VENT PIPE RISER (S)
    ☐_____ MONITORING WELL (S) REQUIRED IN PETROLEUM IMPACTED AREA (S) DESCRIBED IN ITEM 18
    ☐COMPLETE AN ENVIRONMENTAL ASSESSMENT IN COMPLIANCE WITH COMAR 26.10._____(submit two copies)
    ☒SUBMIT SOIL ANALYTICAL RESULTS:
    **EPA METHOD:** ☒8015B GRO/DRO ☒8270 (SVOCs) ☒8260 (VOCs)
    ☒OTHER:PCBS EPA 8081/8082
    ☐SUBMIT GROUNDWATER ANALYTICAL RESULTS:
    **EPA METHOD** ☐8015B GRO/DRO ☐8270 (SVOCs) ☐8260 (VOCs) ☐524.2 (VOCs)
    ☐OTHER _____
    ☒SUBMIT SOIL DISPOSAL RECEIPT
    ☒SUBMIT TANK DISPOSAL RECEIPT

Date: 11/4/15
Case No.: 16-0084BA
Facility No.: 13189

14. (continued)
☒AMEND REGISTRATION: ☐Notification form provided to contact person
☒Owner/Representative informed case file may remain open until notification form is received by MDE
☐ Completed on site.

☐OTHER: ____

15. Has inspector completed: site sketch? ☐YES ☒NO site photographs? ☒YES ☐NO
16. Were tank(s) labeled? (If YES, describe: tank number) ☒YES ☐NO
17. Is follow-up required by this Administration? ☒YES ☐NO

18.COMMENTS:

Writer onsite to observe the removal of seven (7) 1000 gallon steel tanks in Block F portion of LockheEd Martin. The tanks were previously filled in with slurry (concrete/fly ash mix). Drop tubes are inside the tanks and filled with water. Tank 3 and Tank 1 appeared to be partially filled with slurry material and water was observed in tank 3, but no LPH. No perforations were observed in any of the tanks. Piping was encountered on top of the tanks (not connected to the tanks). Piping run is observed at the south corner of tank field leading to the waters edge. The piping run at the tank field end was filled with concrete. The piping will be uncovered and removed at a later date. No water or product was encountered.

Prior to removing the tanks, sheet piling was installed. The tank field was inside a concrete vault with concrete between the tanks. A concrete pad is present at the bottom of the tanks, approximately 12-ft bgs and removed for soil sampling. MAX PID of the soil above and around the tanks were measured at 3-4 ppm ; and were removed for proper offsite disposal. Max PID below the concrete pad was 0.1 - 0.2 units.

REQUIREMENTS:REQUIREMENTS:

1. As per the approved work plan (ROO Work Plan Approval dated 10/6/15) 14 confirmatory soil samples will be collected and analyzed as discussed on site (5 samples will be collected at each side of the excavation (long wall) and below the sheeting if possible. Four samples will be collected within the excavated area, below the tanks).
2. MDE will be notified when piping is exposed for removal and sampling. Soil samples as per the approved work plan will be collected every 15 linear ft.
3. Tanks will be cut open and all the material inside the tanks must be properly disposed of. Tanks will be cut on site and properly disposed. Additional sampling may be required by the disposal facility for waste characterization if the fill is fly ash.
4. Amend the UST Registration.
5. UST Closure Report must be submitted within 45-days including all disposal receipts, laboratory data, site map showing locations of soil samples, tank location, adjacent structures/features, and other information as per the guidance document.

**Ellen Jackson**
Inspector's Name

Ellen Jackson
Inspector's Signature

James R. Ferguson P.E.
Contact Person's Name (printed)

Agent for Lockheed Martin
Contact Person's Signature

410-456-9383
Contact Person's Telephone No.

Larry Macher
Contractor's Name (Printed)

Contractor's Signature

443-324-1103
Contractor's Telephone No.

Lawrence Macher
Technician/Remover Name (printed)

MDIC-2015-1907(R)
Certification Number

May 1, 2017
Expiration Date