# Storm Drainage System Sediment-Sampling Report Lockheed Martin Middle River Complex 2323 Eastern Boulevard Middle River, Maryland

Prepared for:

**Lockheed Martin Corporation** 

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#### **ACRONYMS AND ABBREVIATIONS**

BaPEq benzo(a)pyrene equivalent COC chemical(s) of concern

EESH energy, environment, safety, and health (Lockheed Martin Corporation)

EGIS electronic geographic information system

EROP Enterprise Operation
GC gas chromatography
HASP health and safety plan

IDW investigation-derived waste IRM interim remedial measure

LMCPI Lockheed Martin Corporation Properties, Inc.

Lockheed Martin Corporation

MAA Maryland Aviation Administration

MDE Maryland Department of the Environment

mg/kg milligram(s) per kilogram

µg/kg microgram(s) per kilogram

MRC Middle River Complex

MSA Martin State Airport

NPDES National Pollutant Discharge Elimination System

PAHs polycyclic aromatic hydrocarbons

PCBs polychlorinated biphenyls
PDF portable document format

PM project manager

PPE personal protective equipment
PRG preliminary remedial goal

QA quality assurance
QC quality control

RAL remedial action level

REC recognized environmental condition

SIM selective-ion monitoring

Tetra Tech, Inc.

UCC utility cross-connection

USEPA United States Environmental Protection Agency

VOC(s) volatile organic compound(s)

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## Section 1 Introduction

On behalf of Lockheed Martin Corporation (Lockheed Martin), Tetra Tech, Inc. (Tetra Tech) has prepared this report describing the sampling and chemical analyses of sediments in storm drainage systems that discharge to two waterways adjacent to the Middle River Complex (MRC) and Martin State Airport (MSA) in Middle River, Maryland (see Figure 1-1). The sampling and chemical analyses provide additional sediment chemical data to support the design of the planned sediment remediation in waterways adjacent to the Middle River Complex. These analyses aim to identify sediment in storm drain systems that may act as continuing sources of polychlorinated biphenyls (PCBs) and/or polycyclic aromatic hydrocarbons (PAHs), possibly resulting in PCB concentrations that exceed the preliminary remedial goals (PRGs) and proposed remedial action levels (RALs) that have been established for sediment in Cow Pen Creek and Dark Head Cove (Tetra Tech, Inc., 2013c). This effort also satisfies a request by the Maryland Department of the Environment (MDE) to sample storm drain systems to determine whether sediment in storm drain systems along Wilson Point Road could contribute polycyclic aromatic hydrocarbons (PAHs) to Dark Head Cove sediment (Maryland Department of the Environment, 2013). The storm drainage systems sampled drain both the Middle River Complex (nine systems) and the adjacent Martin State Airport (four systems).

This report supplements the 2013 *Draft Sediment Remedy Design Investigation Report* (Tetra Tech, Inc., 2014a) for the Middle River Complex. Results from this report will be evaluated along with existing sediment data collected during previous sediment investigations (2005–2012), as well as the sediment work completed in 2013 by Tetra Tech, Inc. for Cow Pen Creek and Dark Head Cove. This report is organized as follows:

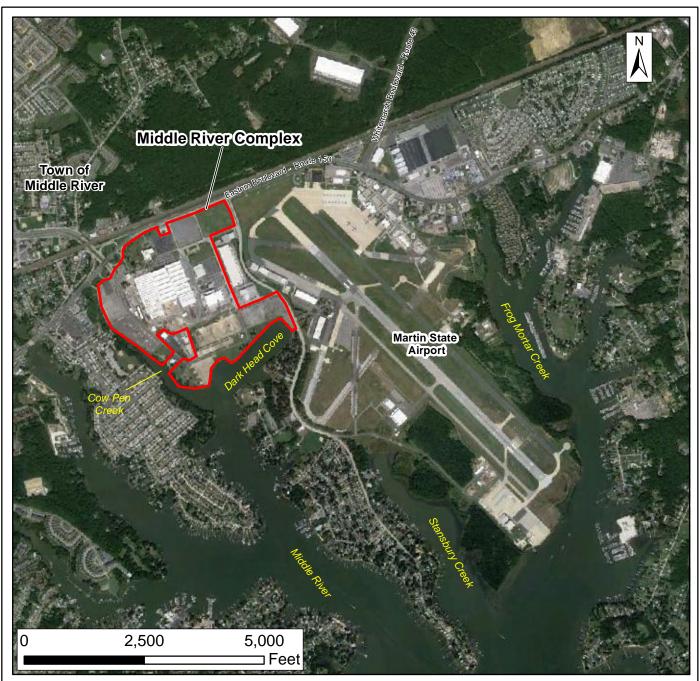
<u>Section 2 – Site Background</u>: Briefly describes the site history and previous investigations related to storm drain sediments at the Middle River Complex.

<u>Section 3 – Investigation Approach and Methodology</u>: Presents the technical approach to the investigation and describes the field methodology for sampling and chemical analyses.

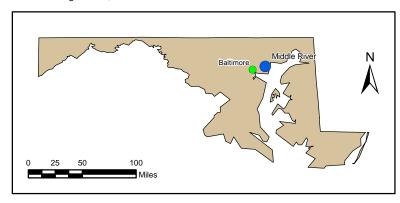
<u>Section 4 – Results</u>: Presents the investigation results.

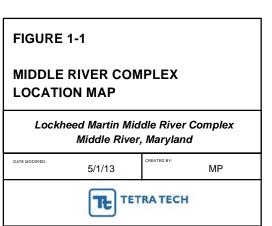
<u>Section 5 – Summarizes</u>: Summarizes the investigation program.

<u>Section 6 – References</u>: Cites references used in compiling this document.



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## Section 2 Site Background

#### 2.1 SITE DESCRIPTIONS

The following sections briefly describe the Lockheed Martin Corporation (Lockheed Martin) Middle River Complex (MRC) and Martin State Airport (MSA). Details of the area's physical setting, land use, physiography, and subsurface conditions (i.e., soils and geology), and previous studies related to creek-bottom sediments in Cow Pen Creek and Dark Head Cove, are in the 2013 *Draft Sediment Remedy Design Investigation Report* (Tetra Tech, Inc. [Tetra Tech], 2014a), and are not repeated herein.

#### 2.1.1 Middle River Complex

The MRC is at 2323 Eastern Boulevard in Middle River, Maryland. Figure 2-1 is a facility layout map. The site is comprised of approximately 163 acres and 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, MSA to the east, Dark Head Cove to the south, and Cow Pen Creek to the west. As shown in Figure 2-1, the MRC currently consists of eight tax parcels or "tax blocks," which are commonly referred to as Block A, B, D, etc.

Currently, the primary activities of LMC Properties, Inc. (LMCPI) (a subsidiary of Lockheed Martin) at the MRC include facility and building management and maintenance. The main site tenant, MRA Systems, Inc. (a subsidiary of General Electric Company), designs, manufactures, fabricates, tests, overhauls, repairs, and maintains aeronautical structures, parts, and components for military and commercial applications. Lockheed Martin Mission Systems & Training is a business segment of Lockheed Martin that conducts engineering activities and fabricates, assembles, tests, and otherwise supports vertical-launch systems. A subsidiary of Lockheed

Martin (Applied NanoStructured Solutions LLC) occupies a portion of the MRC, where it researches and designs nanotechnology applications.

Environmental investigations of the MRC were substantially initiated with a 2003 environmental site assessment which identified several recognized environmental conditions associated with the MRC (Earth Tech, Inc., 2003). Since that time, extensive environmental investigations have been conducted at the MRC and remedies are being implemented under several remedial action plans to mitigate conditions in soil, groundwater, and sediment under a consent order to be issued by the Maryland Department of the Environment (MDE).

#### 2.1.2 Martin State Airport

MSA is at 701 Wilson Point Road in Middle River, Maryland. It is bounded by Frog Mortar Creek to the east and Stansbury Creek to the west (see Figure 2-1). Both creeks are tidal tributaries of the Chesapeake Bay that intersect Middle River, which flows to the bay south of the airport. The current MSA property was part of the Glenn L. Martin Company's original 1,260-acre property, which previously included a manufacturing portion (present-day MRC), as well as previously divested properties now occupied by the Tilley Chemical Company, North American Electric Company, an Exxon service station, the United States Postal Service, the Annex building (currently leased by Ace Logistics, Inc.), and Johnson & Towers, Inc. On September 20, 1975, the Maryland Aviation Administration (MAA) purchased 747 acres that are now used as the airfield, of which 175 acres in the northeastern portion are now leased to the Maryland Air National Guard.

MAA currently operates MSA on behalf of the Maryland Department of Transportation. MSA has an administration building (main terminal building), aircraft hangars, a 7,000-foot long runway, several taxiways, and the Strawberry Point Hangar. MAA manages more than 130,000 square feet of heated hangar space and 190 smaller aircraft T-hangars. The southwestern portion of MSA contains numerous aboveground fuel storage tanks for Jet A and Avgas 100LL fuels. MSA is also home to more than 20 commercial tenants providing fuels and lubricants, helicopter avionics repair, and flight instruction, in addition to hosting the Baltimore County Police aviation and marine units and the Glenn L. Martin Museum (MAA, 2012). Lockheed Martin currently conducts a variety of environmental investigations at MSA, under MDE review and regulation.

#### 2.2 STORM DRAIN SYSTEMS

#### 2.2.1 Middle River Complex

No surface water bodies lie within or cross the MRC. Excluding areas immediately adjacent to Cow Pen Creek and Dark Head Creek, surface water runoff from the MRC discharges from the facility via storm drains. In the outer tax blocks, runoff may also enter creeks directly via overland flow from paved or concrete surfaces (e.g., the sloped concrete ramp in Block F), or in vegetated areas near the creek by overland flow, rills, or gullies. Water from runoff may also infiltrate into the soil and move as shallow subsurface storm-flow, or may locally pond and infiltrate into the soil or be lost to the atmosphere through evaporation.

Stormwater in the industrial portions of the MRC is collected and conveyed through nine primary storm-drainage systems that discharge to Cow Pen Creek and Dark Head Cove via Outfalls 001 through 009 (Figure 2-2). Discharges from These outfalls are regulated under a State of Maryland National Pollutant Discharge Elimination System (NPDES) permit (State discharge permit No. 00-DP-0298, NPDES permit No. MD0002852) issued by the MDE Industrial Discharge Permits Division, Water Management Administration). Four outfalls are monitored monthly for flow, chlorine and pH (Outfalls 002, 004, 007 and 009); Outfall 004 is monitored monthly for oil and grease; and Outfalls 004 and 009 are monitored quarterly for total and dissolved copper. Additionally, stormwater in portions of Block G and Block H is collected and conveyed to Cow Pen Creek by other smaller storm-drainage systems that are not covered under the permit.

#### 2.2.2 Martin State Airport

Surface water runoff from MSA enters surrounding creeks via localized gullies in the eastern and western undeveloped portions of the site, or via storm drains that collect and convey runoff from the airport runway, taxiways, and developed administrative portions of the facility. MSA encompasses 47 drainage areas in three watersheds, forming a total drainage area of 700 acres (MAA, 2008). The airport drainage areas range from seven to more than 170 acres. The western MSA boundary consists of the main terminal building and aircraft hangars. Stormwater runoff from the eastern and southern ends of Dark Head Cove flows via five storm drain systems to MSA outfalls that are monitored for water quality (90F001, 80F001, IN159, SW0043, and WROF001A). The monitored outfalls are on MSA property (Figure 2-3), but stormwater from

MSA is conveyed to Dark Head Cove by storm drains running beneath Wilson Point Road, MRC Block D, and Stansbury Apartments.

All stormwater runoff originating from MSA discharges to outfall areas that are monitored monthly to ensure that no oily discharges to surface water occur. Secondary containment drains are also routinely inspected and emptied of stormwater. The facility maintains an NPDES permit (No. MDR 05501, General Discharge Permit No. 05-SF-5501), with an effective date of November 12, 2004, and an expiration date of November 12, 2009. However, the current permit remains in effect because the MDE has administratively extended it until they issue a replacement.

MSA's general industrial permit has no monitoring requirements. However, limited monitoring is performed for the separate municipal permit required for the federal Illicit Discharge Detection and Elimination program. This limited monitoring includes laboratory analysis for ammonia, dissolved oxygen, surfactants, fecal coliform, potassium, water temperature, conductivity, pH, and fluoride concentrations in monitored outfalls during annual inspections. Visual inspections are routinely performed and annual reports are submitted to MDE.

#### 2.3 PREVIOUS INVESTIGATIONS

The following sections summarize previous environmental investigations and an interim remedial measure (IRM) implemented for the chemical characterization and removal of storm drain sediments at the MRC. Lockheed Martin has not conducted previous studies of storm drain sediment at the MSA. However, Lockheed Martin has conducted studies of sediment in Dark Head Cove that include sampling downstream of several MSA storm drain outfalls. Detailed summaries of Dark Head Cove sediment studies are in the 2013 *Draft Sediment Remedy Design Investigation Report* (Tetra Tech, 2014a).

#### 2.3.1 Supplemental Soil and Storm-Drain Sediment Characterization Investigation (Summer 2009)

In 2009, Lockheed Martin investigated sediment in Block E storm-drain systems that discharge to Dark Head Cove (Tetra Tech, 2010). Sediment samples were collected from storm drains to determine if Block E chemicals of concern (COC) (i.e., PCBs, polycyclic aromatic hydrocarbons [PAHs], volatile organic compounds [VOCs], and metals) were present in storm drain sediments,

indicating that these sediments may be past or continuing sources of COC to bottom sediments of Dark Head Cove. Sediment samples were collected from 20 storm-drain catch basins (i.e., grated inlets) and manholes within three storm-drain systems associated with MRC Outfalls 005, 006, and 008. Samples were analyzed at an off-site laboratory for PCBs, PAHs, VOCs, and metals.

PCBs were detected in nearly all sediment samples collected from Block E storm drains, with maximum PCB concentrations (downstream to upstream) of 91 milligrams per kilogram (mg/kg), 102 mg/kg, and 31 mg/kg. All total PCB concentrations in sediment from the Outfall 005 system exceeded the preliminary sediment-cleanup level used at the time of the study (one milligram per kilogram [mg/kg]), and all but one concentration exceeded 10 mg/kg. PAH concentrations (evaluated as benzo(a)pyrene equivalents [BaPEq]), 0.332-139 mg/kg, with all but three sample concentrations exceeding the preliminary BaPEq cleanup-goal for sediment that was used at the time (0.70 mg/kg). Cadmium concentrations in three samples, and the chromium concentration in one sample, also exceeded the preliminary cleanup goals (eight and 138 mg/kg, respectively) used at the time of the study. VOCs were detected only at trace to low concentrations. Therefore, the project team concluded that PCBs, PAHs, and metals concentrations in Dark Head Cove sediments may be partially attributable to sediment discharging from these three outfalls.

#### 2.3.2 Utility Cross-Connection Study (Summer 2011)

A utility cross-connection (UCC) investigation (Tetra Tech, 2012d) was conducted in 2011 of areas in Blocks G, I, and E that had been identified for *in situ* groundwater remediation in the final *Groundwater Response Action Plan* (Tetra Tech, 2012b). The UCC investigation included the southeastern portion of Block G, the area immediately south and southeast of Building C in Block I, and the southeastern portion of Block E near a 500,000-gallon water tank and REC #3. The investigation sought to identify utilities within the footprint of the proposed groundwater-remedy systems that may act as preferential migration pathways for groundwater contaminants or bioremediation substrates that may be injected into groundwater to remedy VOCs. The UCC investigation results were used to design pilot-scale groundwater-injection tracer tests (Tetra Tech, 2012c) and the full-scale groundwater remedial system designs Tetra Tech, 2013d). No environmental samples were collected for subsurface characterization during the UCC study.

The UCC investigation included reviews of historical drawings, employee interviews, site reconnaissance, geophysical surveys, visual and closed-circuit television inspections of storm drains, and professional land surveying to locate, record, and map subsurface utilities. Historical and current utility records, site reconnaissance, and geophysical surveys indicated numerous underground utilities in the three study areas. Underground utilities include electrical lines, telecommunication lines, domestic- and fire-water lines, sanitary sewer lines, natural gas lines, and storm drains. Known utilities and geophysical anomalies were marked in the field. These locations were professionally surveyed to provide location coordinates and elevations, and to provide detailed maps of the utilities for future remedial designs.

Closed-circuit television was used to inspect and digitally record accessible underground structures. The storm drains were cleaned to remove sediment and debris before the closed-circuit television inspection, to allow for free movement of the crawler camera. Cleaning involved inserting a jet nozzle into the downstream structure of the storm drain and propelling it toward an upstream structure.

As the nozzle was pulled back, liquid and solid material was vacuumed into a jet-vacuum truck. Jet-vacuum-truck contents were transferred to roll-offs or fractionation tanks (located on a containment pad) for characterization and disposal. After a pipe segment had been cleaned, the mobile closed-circuit television truck was positioned at the upstream structure and a robotic crawler camera equipped with a multi-angle lens was inserted into the drainage pipe for video inspection of the downstream structure. All observations were recorded on a hard drive, and an audio commentary accompanied the video inspection.

The results of the utility cross-connection study indicate that the storm-drain and sanitary-sewer lines are at similar elevations to the groundwater table in the three areas where injections of Lactoil<sup>®</sup> into groundwater are proposed. Block G and Block I were historically active areas, and contain used and abandoned sewers and storm drains. Block G injection areas once contained multiple buildings and a wastewater treatment plant.

Proposed injection areas in Blocks E, G, and I also contain geophysical anomalies that might or might not affect distribution of Lactoil<sup>®</sup>. Groundwater was observed flowing in a conduit joint in the Block E injection area. Previous pilot work and the results of the utility cross-connection investigation show the potential for injected Lactoil<sup>®</sup> substrate to enter the storm drain system or

other utility conduit resulting in a possible release to the nearby surface water body. The report recommended pilot-scale bromide-injection tracer tests for the three proposed injection areas to determine optimal substrate-injection rates and evaluate whether injected substrate and VOCs in groundwater would enter the storm-drain or sanitary-sewer systems. Low-flow injection rates were recommended for full-scale implementation, to limit groundwater mounding in the areas where utilities and groundwater were at similar elevations.

#### 2.3.3 Block E Storm-Drain Interim Remedial Measure (Fall 2011)

A storm-drain IRM was completed at Block E in autumn 2011 (Tetra Tech, 2012a). The IRM removed sediment and debris from Block E drainage-system piping and manholes, provided sediment controls, and repaired or replaced inlets and manholes. The IRM was intended to minimize the transport of PCB-contaminated storm-drain sediments to off-site locations (e.g., Dark Head Cove) and to identify and remove flow restrictions where possible. The IRM was not designed to achieve preliminary sediment-cleanup goals, but rather to remove the more mobile sediment that might migrate off-site through the drainage system. The IRM was an initial step in the more comprehensive remediation of PCB-contaminated sediments planned for the Dark Head Cove area. No environmental sampling was conducted as part of the storm-drain IRM except during management of waste sediment removed from the storm drains (i.e., waste profile sampling and analyses).

A Maryland-licensed land surveyor conducted a pre-construction survey to record topographic and as-built features of the planned and potential disturbance areas around the Block E storm-drain system. Following completion of all fieldwork, a post-construction survey documented that the site had been restored to a condition as close as practical to its original pre-construction state. Inlets, manholes, or pipe sections of the storm drain system that were found to be damaged beyond repair or blocked and inaccessible at the beginning of the IRM were generally left as found. The accessible storm drain lines were cleaned of mobile sediment.

Cleaning the storm drains involved inserting a jet nozzle into the downstream structure and propelling it toward the upstream structure. As the nozzle was pulled back, liquid and solid material was vacuumed into a jet-vacuum truck. Hand removal of debris was necessary at some locations due to the large size or heavy weight of the debris in these structures. Jet-vacuum-truck contents were transferred to roll offs or fractionation tanks located on a containment pad. After a

pipe segment had been cleaned, the mobile closed-circuit television truck was positioned at the upstream access structure, and a robotic-crawler camera equipped with a multi-angle lens was inserted into the drainage pipe to video inspect the downstream structure. All observations were recorded on a hard drive, and an audio commentary accompanied the video inspection.

Three additional catch basins were identified in the northeast corner of Block E during a site walk in July 2011, before implementation of the Block E storm-drain system IRM. Two buried manholes (MH-8 and MH-9) were also found near the southeast corner of the former Building D foundation during the IRM. These manholes have grated-steel covers and previously served as catch basins for local runoff. Manhole MH-9 was raised to grade, and a new cover was placed over the manhole opening. Manhole MH-9 is filled with rocks that partially obstruct flow from the upstream pipe; however, some water flows through manhole MH-9 to the downstream catch basin and piping. MH-8 could not be cleaned or filmed because piping was filled with water caused by rocks obstructing the connection from MH-9.

Repairs using brick, concrete, grout, and cast-iron frames and covers were performed as the cleaning of the Block E storm-drain system progressed. Crushed stone backfill and concrete or vegetative surface finishes were used as needed to return the structure/location to its original or better condition. Upon completion of the IRM, all disturbed areas were restored to approximately pre-existing grades. Seed and topsoil were distributed across all disturbed vegetated areas. Land surveys, photographs, and a post-cleaning video survey of the Block E drainage pipes document site restoration activities. Sediment controls, including silt fences and hay bales, were then installed around catch basins and manholes, so that sediment reentering cleaned drain lines was minimized until the area could be stabilized with grass.

#### 2.4 PROBLEM STATEMENT AND CURRENT INVESTIGATION

Various MRC site investigations from 2003–2012 have identified sediment contamination in Cow Pen Creek and Dark Head Cove resulting from historical facility activities (Tetra Tech, 2014a). Sediment impacts include elevated concentrations of PCBs, PAHs, and metals. Sediment sampling results indicate possibly unacceptable (according to environmental regulations) risks to local fauna, including benthic (i.e., sediment dwelling) invertebrates, fish, and piscivorous (fish-eating) birds. Risk drivers include PCBs and, to a lesser extent, PAHs, present in the top 30 inches of the sediment bed. Specifically, PCBs and PAHs were identified in shallow

sediments near the bulkhead; however, only PAHs were identified in Dark Head Cove near MSA. In addition, cadmium and chromium (metals potentially toxic to benthic macroinvertebrates) were found in sediment at deeper intervals. Both metals have also been identified in Cow Pen Creek and Dark Head Cove sediments.

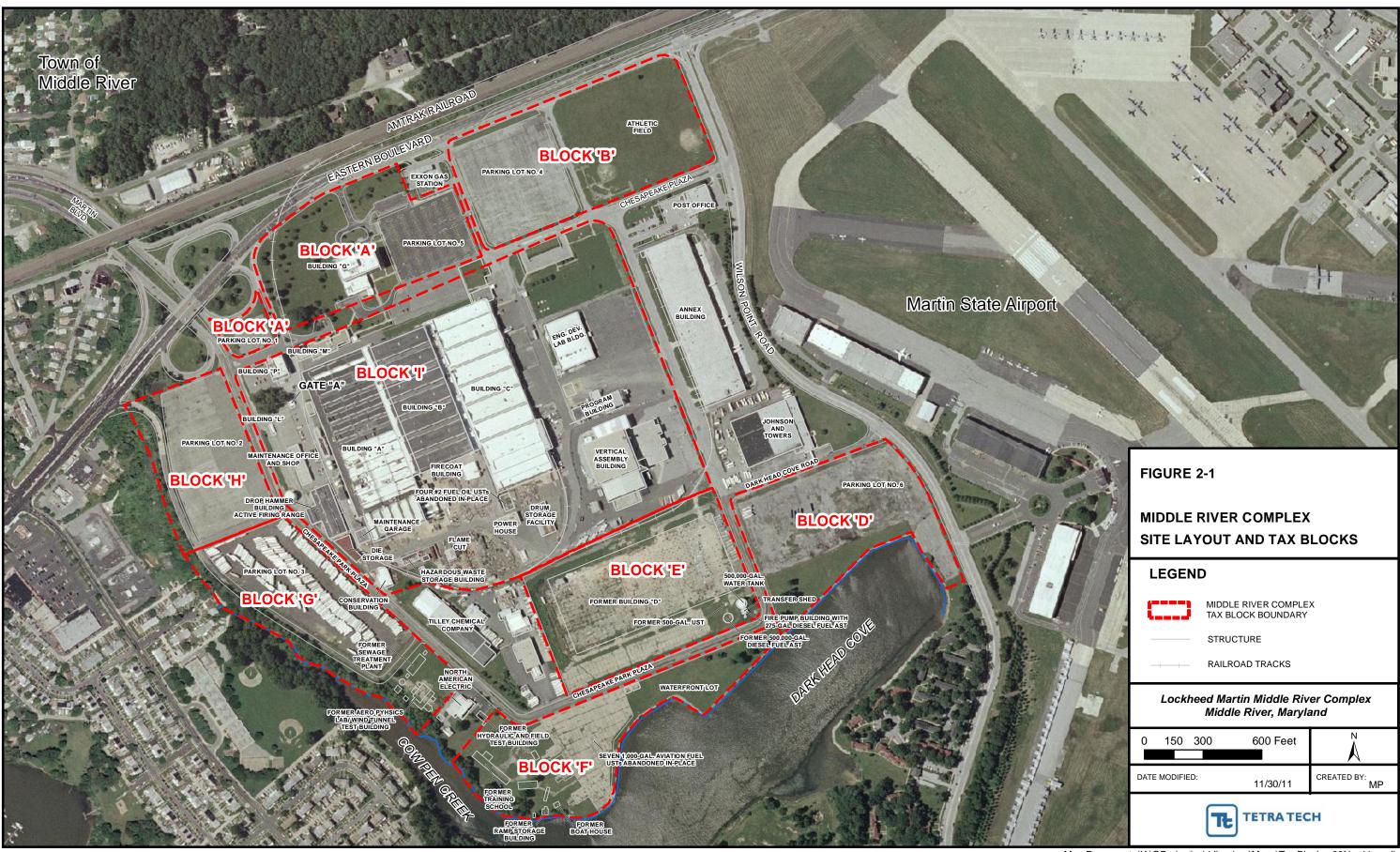
Currently, remediation of sediment is planned for areas where PCBs, PAHs, and metals exceed both proposed cleanup goals (known as preliminary remediation goals [PRGs]) and removal/treatment concentrations (known as remedial action levels [RALs]). RALs identify areas that require treatment or removal to achieve the risk-based PRGs. Site sediments are also being studied to assess the suitability of treating contaminated sediments *in situ* (i.e., in place) to reduce COC bioavailability and minimize disturbance of the habitat.

Other MRC studies have investigated the occurrence and distribution of PCBs, PAHs, VOCs, and metals in surface and subsurface soils in areas adjacent to Cow Pen Creek and Dark Head Cove. Detailed summaries of these soil investigations are in the Tax Blocks E and F soil remedial action plans (Tetra Tech, 2013b and 2014b). At two land parcels (i.e., Tax Blocks E and F) near Dark Head Cove, PAHs were detected in soil at concentrations associated with regulatorily unacceptable human health risks for hypothetical industrial receptors. PCB concentrations in Block E also exceeded acceptable levels in soil (50 mg/kg) according to regulations under the Toxic Substances Control Act of 1976 (40 Code of Federal Regulation 761).

Previous sediment sampling results indicate that sediment in Dark Head Cove has been impacted by contamination transported to the cove via storm drains. For example, elevated PCB concentrations have been detected in surface soil near Block E catch basins and in Block E storm-drain sediment (before the storm drain IRM), and elevated PCB concentrations are found in Dark Head Cove sediment samples collected adjacent to the Block E storm-drain outfalls. Sediment in storm drain systems discharging to Cow Pen Creek (Outfalls 001 through 004) have not previously been sampled and chemically characterized for PCBs, PAHs, and metals.

Additionally, the Block E storm drain systems have been active for nearly two years since the IRM was completed in autumn 2011. Therefore, new chemical data are required for active storm-drain systems discharging into Cow Pen Creek and Dark Head Cove, to evaluate if sediment in these systems could, following remediation, act as continuing sources of PCB and/or PAH contamination to sediment in these two water bodies. Sediment in storm drain utilities was

sampled at both MRC and MSA; this report details the scope-of-work performed to address the investigation objective. This work also satisfies a request by MDE to sample storm drain systems to determine whether sediment in storm drain systems along Wilson Point Road would continue to contribute PAHs to sediment of Dark Head Cove (MDE, 2013).





**400 FEET** 

OUTFALL

STORM WATER INFILTRATION TRENCH

MIDDLE RIVER COMPLEX AND MARTIN STATE AIRPORT, MIDDLE RIVER, MARYLAND

BASED ON TAI 2002 UTILITY PLAN, MARTIN STATE AIRPORT UTILITY COVERAGE, AERIAL PHOTOGRAPHS, TETRA TECH FIELD SURVEYS, AND TETRA TECH SITE INSPECTIONS; ALL UTILITIES BASED ON TAI 2002 UTILITY PLAN AND MARTIN STATE AIRPORT COVERAGE NOT VERIFIED IN THE FIELD

## Section 3 Investigation Approach and Methodology

Tetra Tech, Inc. (Tetra Tech) collected samples of accumulated sediment for laboratory chemical analyses from nine Middle River Complex (MRC) and four Martin State Airport (MSA) storm drainage systems that discharge to Cow Pen Creek or Dark Head Cove (Figure 3-1). Tetra Tech obtained permission from MSA before sampling began. This investigation entailed the following field and post-field activities:

- mobilization/demobilization of sampling staff and equipment
- collection of sediment samples from storm drainage systems that discharge into Cow Pen Creek and Dark Head Cove
- collection, storage, and characterization of investigation-derived waste (IDW), and disposal of that waste at an off-site Lockheed Martin-approved disposal facility
- laboratory chemical analyses and chemical-data validation on the sediment samples
- evaluation of the sampling results

#### 3.1 MOBILIZATION/DEMOBILIZATION

Following regulatory approval of the work plan (Tetra Tech, 2013a), the Tetra Tech field operations leader coordinated mobilization, which included procuring the required subcontractor (laboratory) and mobilizing personnel and materials to the field. Mobilization also included locating the appropriate equipment required for the field tasks, purchasing necessary equipment as required, and staging equipment for efficient loading and transportation to the site. Mobilization began on December 17, 2013 and included the following:

- mobilization of equipment and materials to the site
- implementation of:
  - o site-specific health and safety plan (HASP)

- o emergency response plan
- o sampling and analysis plan
- o waste management plan conforming to Lockheed Martin's *Energy, Environment, Safety, and Health (EESH) Remediation Waste Management Procedure No: EROP-03, Revision 4* (effective April 17, 2009) (Lockheed Martin. 2009a)
- o quality assurance (QA)/quality control (QC) plan
- o data management plan
- set up a decontamination area

#### Demobilization included:

- demobilization of equipment and materials from the site
- general cleanup and trash removal
- management of IDW

Field activities planned for the locations identified in this investigation were coordinated with Lockheed Martin, LMC Properties, Inc. and Martin State Airport (MSA). No drilling or excavating was performed for this study; therefore, no utility clearance was required before sampling.

#### 3.2 SEDIMENT SAMPLING AND ANALYSES

Storm-drain-sediment sampling locations are shown in Figure 3-1. Sediment samples were collected December 18–20, 2013 from the nine primary MRC storm-drainage systems (Outfalls 001–009) and four MSA storm-drain systems (Outfalls 90F001, 80F001, IN159, and WROF001A) that discharge to Dark Head Cove. Attempts were made to collect sediment samples from accessible locations (e.g., manholes or catch basins) containing sufficient sediment to obtain a sample aliquot upstream of the discharge point (i.e., outfall) for each storm drain system. Many access points lacked sediment, or lacked sufficient sediment for the required sample aliquot. Therefore, manholes or catch basins sampled for several systems were located substantially upstream of outfalls (e.g., MRC Outfalls 001, 002, 006, and 007), or in some cases only one or two samples per system could be sampled (e.g., MRC Outfalls 003 and 004, and MSA Outfalls 80F001 and IN159).

Thirty-two sediment samples were collected, with sequential sample designations SD-163 to SD-194. SW0043 is an MSA outfall to Dark Head Cove from an infiltration basin that receives direct runoff from a nearby building, grassy areas, and a taxiway. The basin has no upstream catch basins or piping. Sediment was not present in the infiltration basin or in the outfall manhole near the property boundary; therefore, a sediment sample could not be collected for this stormwater system.

Grab samples were collected from each location using a flat, stainless steel scoop (AMS, Inc. Model 428.06 #12). The sample aliquot was homogenized using the scoop and stainless steel bowls. After each sample had been homogenized, the sediment was placed into a sample container supplied by the analytical laboratory, and the sample container was then immediately placed in a cooler containing ice. All reusable equipment contacting sediments was decontaminated between sampling locations, as described in Section 3.4. Sediment samples were analyzed for polychlorinated biphenyls (PCBs), polycyclic aromatic hydrocarbons (PAHs), and percent moisture using methods listed in Table 3-1.

Sample descriptions were recorded, noting color, grain size, sorting, texture, odor, and any other pertinent sediment characteristics. Grain size was determined by comparing sediment grains to a grain-size chart. Location and construction information for each sampled catch basin or manhole were recorded. Location information includes the facility location (MRC or MSA), outfall number, catch basin or manhole number (if available), and a location sketch. Construction information includes material type (i.e., concrete, brick and mortar, etc.); utility dimensions (length, width, bottom depth); an estimate of the amount of sediment present (i.e., area and thickness); and the number, size, material, and locations of inflow and outflow structures. Inflow/outflow structures could not be observed in flooded catch basins or manholes, and the presence and quantity of sediment could only be estimated by the field crew using sampling tools and probes. The field information was documented on a Tetra Tech sediment-sampling form (Appendix A).

#### 3.3 SAMPLE NOMENCLATURE AND HANDLING

Sediment samples submitted to the laboratory were labeled with an "SD" prefix, identifying the sampled medium as sediment, followed by a three-digit numeral to identify the sampling location (numbers 163–194), followed by the sampling depth-interval, beginning at the top of the

sediment surface. For example, SD-184-0–2 designates a surface sediment sample collected at location SD-184 from a sediment depth-interval of zero to two inches. Proper custody procedures were followed throughout all phases of sample collection and handling. Laboratory-cleaned sample containers were released under signature from the laboratory and were accepted under signature by the samplers. Transport containers were sealed with strapping tape and a tamper-proof custody seal. The custody seal contains the signature of the individual releasing the transport container, along with the date and time.

Chain of custody protocols were used throughout sample handling to establish the evidentiary integrity of sample containers. These protocols demonstrate that the samples were handled and transferred in a manner that would prevent tampering. Sample chain of custody forms are in the data validation reports in Appendix B.

#### 3.4 EQUIPMENT DECONTAMINATION

Reusable sampling equipment (scoop and bowls) was decontaminated between sampling locations before each use. Decontamination solutions were collected for appropriate disposal. Decontamination consisted of the following steps:

- Liquinox® and potable-water wash
- Potable-water rinse
- Reagent-grade isopropanol rinse (achieved by thoroughly wetting the equipment with isopropanol)
- Analyte-free water rinse
- Air drying

#### 3.5 WASTE MANAGEMENT

The project followed a waste management plan conforming to *Lockheed Martin EESH Remediation Waste Management Procedure EROP-03, Revision 4* (Lockheed Martin, 2009a) and *Remediation ESH Contractor Handbook*, (Lockheed Martin, 2009b). IDW, consisting of equipment rinse water and personal protective equipment (PPE), was generated during this sampling event. PPE was placed in trash bags and disposed of in a Lockheed Martin-designated trash container. All drums were appropriately labeled and logged on a drum inventory form. IDW

was characterized and disposed of in accordance with applicable state and federal regulations. IDW generated during this investigation was disposed of as non-hazardous waste.

#### 3.6 DATA MANAGEMENT

Data handling procedures followed by the laboratory met the requirements set forth in the laboratory subcontract. All analytical and field data are maintained in Tetra Tech project files. The project files contain copies of the chain of custody forms, sampling log forms, sampling location maps, and QA/QC documentation.

#### 3.6.1 Data Tracking and Control

A cradle-to-grave sample-tracking system was used throughout the investigation. Before field mobilization, the field operations leader coordinated and initiated sample tracking. Sample labels were preprinted before entering the field with sample ID, date, and time, then handwritten in the field. Labels were reviewed for accuracy and for adherence to work plan requirements. The project manager (PM) coordinated with the analytical laboratory personnel to ensure that they were aware of the number and type of samples and analyses they would receive. When field sampling was underway, the field operations leader forwarded the chain of custody forms to the PM or designee and to the laboratory at the end of each day of sampling. The PM/designee confirmed that the chain of custody forms provided the information required by the work plan.

This data management system ensures early detection of possible field errors so adjustments can be made while the field team is mobilized. After successful completion of all requested analyses, the laboratory submitted an electronic deliverable for every sample delivery group. When all electronic deliverables have been received from the laboratory, the PM/designee ensured that the laboratory performed all requested analyses.

#### 3.6.2 Sample Information

Data from field measurements were recorded using the appropriate log sheets (see Appendix A). Reduction of field data entailed summarizing and presenting the data in tabular form. Reducing laboratory data entailed manipulating raw data-instrument output into reportable results. Laboratory data were verified by the group supervisor and by the laboratory's QC documentation department.

#### 3.6.3 Project Data Compilation

The analytical laboratory generated an Adobe *Acrobat*<sup>®</sup> portable document format (PDF) file of the analytical data packages, as well as electronic database deliverables. The electronic database was checked against the PDF file provided by the laboratory and updated based on data-qualifier flags applied during data validation, as required. Sediment data were incorporated into the electronic geographic information system (EGIS) database. All data, such as units of measure and chemical nomenclature, were reviewed and corrected, if necessary, to maintain consistency with the project database.

#### 3.6.4 Geographical Information System

Data management systems now in use consist of a relational database and an EGIS to manage MRC environmental information. The relational database stores chemical, geological, hydrogeological, and other data collected during environmental investigations. The EGIS is built from the relational database and contains subsets of the larger data pool. The EGIS can post environmental data onto base maps to represent the information graphically. Upon compilation of sample, chemical, and positional data, the data were incorporated into the MRC and MSA EGIS.

#### 3.7 DATA VALIDATION

A party independent of the analytical laboratory validated the data by reviewing it to ensure that specific criteria had been met. These criteria are concerned with specifications that are not sample-dependent; they specify performance requirements that should be fully under a laboratory's control. For organic-data analyses, specific validation areas include blanks, performance-evaluation-standard materials, and instrument-performance checks. The chemical data were validated by Tetra Tech's data validation group in Pittsburgh, Pennsylvania.

Chemical data were supplied by the laboratory as hard-copy reports and electronic databases. After sampling was complete, chemical data were validated to assess their reliability and accuracy, in accordance with established United States Environmental Protection Agency (USEPA) protocols. This review was based on USEPA Region 3 guidelines (USEPA, 1993 and 1994) and the specifics of the analytical method employed. Data validation reports with chains of custody and a data table of all sampling results (including non-detects) are in Appendices B (data validation reports) and C (data table).

Collectively, these data are acceptable for their intended use. For this validation, the following data qualifiers (i.e., flags) were applied to chemical results presented in this report:

- J The analyte is considered present in the sample; however, the value is estimated and may not be accurate or precise.
- U Not detected; the analyte is considered not detected at the reported value.

The qualifiers appear in the chemical-results tables in Section 4 and Appendix C, and in the data validation report in Appendix B.

Table 3-1
Sampling and Chemical Analyses for Storm Drainage System Sediment Samples—2013
Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland

Sampling location and number		Sampling depth	Sample analyses and methods	Rationale/purpose
SD-163 through SD-194 (32 sediment samples)  MRC Outfall 001 SD-163-0-1 SD-164-0-1 SD-165-0-1 MRC Outfall 002 SD-166-0-1 SD-168-0-1 SD-168-0-1 MRC Outfall 003 SD-169-0-1 MRC Outfall 004 SD-170-0-1 SD-172-0-3 SD-173-0-3 SD-181-0-1 SD-182-0-3 SD-183-0-1 MRC Outfall 006 SD-187-0-1 SD-188-0-1 SD-189-0-1	MRC Outfall 007 SD-190-0-1 SD-191-0-1 MRC Outfall 009 SD-184-0-2 SD-185-0-1 SD-186-0-1 MSA Outfall 9OF001 SD-176-0-3 SD-177-0-2 SD-180-0-1 MSA Outfall 8OF001 SD-178-0-1 SD-179-0-1 MSA Outfall IN159 SD-174-0-1 SD-175-0-1 MSA Outfall WROF001 (downstream) SD-192-0-1 SD-193-0-1 SD-194-0-1	0–1 inch 0–2 inches or 0–3 inches	Laboratory analyses: PCBs by SW846 Method 8082  PAHs by SW846 Method 8270 (GC/SIM)  percent moisture	Sample and chemically analyze sediments that are currently in storm drainage utility lines that could represent a continuing source of PAHs or PCBs to sediments of Cow Pen Creek and Dark Head Cove.

GC = gas chromatography

MRC = Middle River Complex

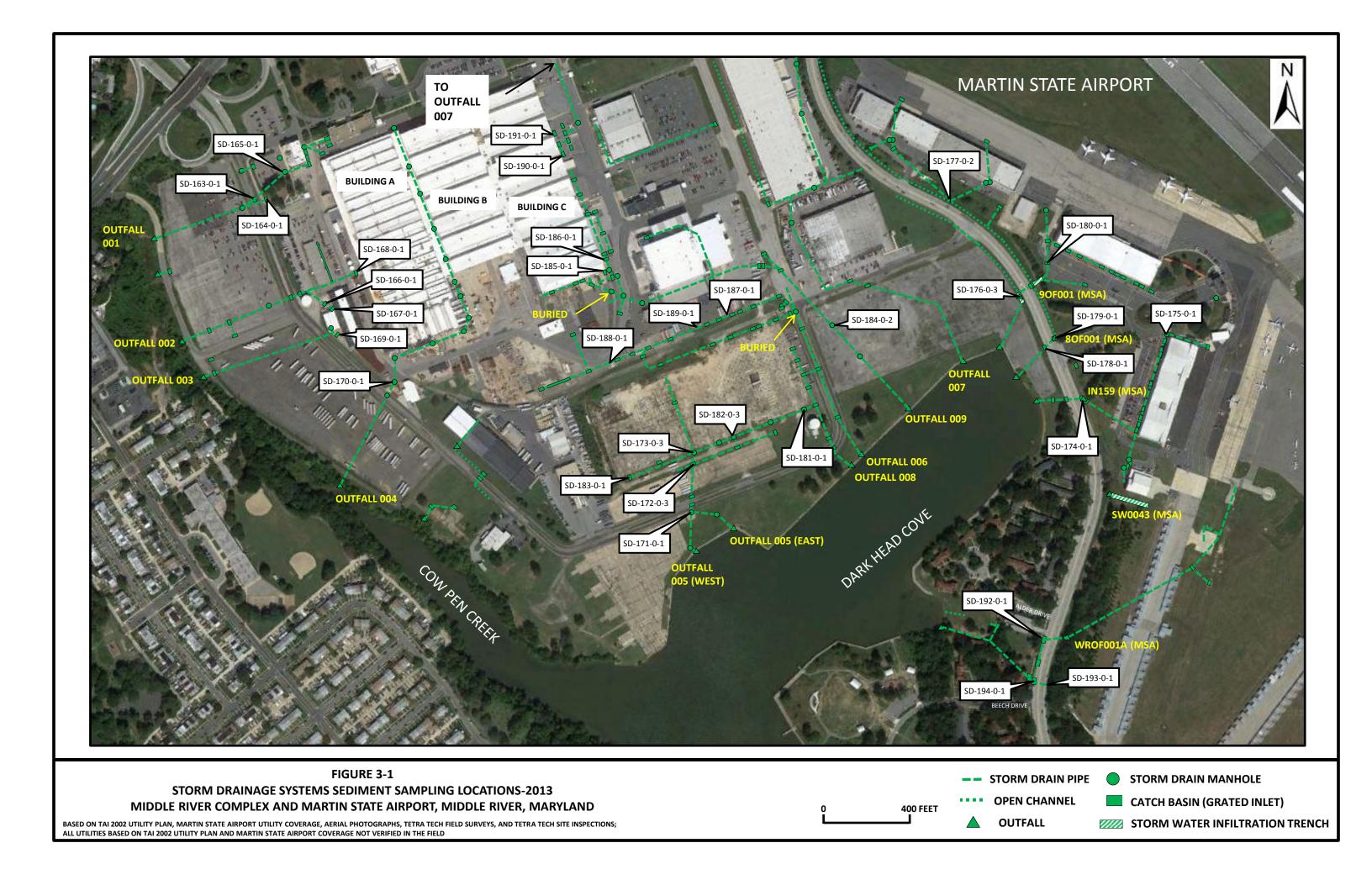
MSA = Martin State Airport

PAHs = polycyclic aromatic hydrocarbons

PCBs = polychlorinated biphenyls

SD = sediment

SIM = selective-ion methodology



### Section 4 Results

This section presents the results of the 2013 storm-drain-sediment field investigation. This investigation sought to provide data to chemically characterize and identify sediments currently in storm-drain utility lines, because they may represent a continuing source of polychlorinated biphenyls (PCBs) or polycyclic aromatic hydrocarbons (PAHs) to Cow Pen Creek or Dark Head Cove sediment. These data were collected to support the remedial design for sediment in Cow Pen Creek and Dark Head Cove.

Most catch basins and manholes inspected lacked sediment or lacked sufficient sediment for a sample aliquot. Most appurtenances sampled contained one inch or less of sediment. These observations were not unexpected, because prior camera studies of the storm drain system (Tetra Tech, Inc. [Tetra Tech], 2012a, 2012d) have shown that sediment tends to accumulate at pipe joints (particularly offset joints or mineral encrustations), and not necessarily in accessible portions of catch basins and manholes. One manhole at sampling location SD-173-0–3 (at a manhole that could not be cleaned as part of the 2011 storm drain interim remedial measure [IRM]) was entirely filled with soil.

Validated chemical data for the 32 storm-drain-sediment samples were used to generate a statistical summary table (Table 4-1) and a table summarizing positive detections of the chemical analytes (Table 4-2). Samples were analyzed for PCBs and PAHs. Total PCBs and benzo(a)pyrene-equivalent (BaPEq) concentrations were calculated for each sample and these were compared to the proposed point-based (i.e., single sample) remedial action levels (RALs) in the Middle River Complex (MRC) sediment feasibility study (Tetra Tech, 2013c). PAH concentrations for each sample were typically converted to a BaPEq concentration so that risk could be evaluated using a single risk-based value for PAHs. The BaPEq is the sum of all carcinogenic PAH concentrations that have been adjusted by applying chemical-specific toxicity-equivalence factors relative to benzo(a)pyrene. For reference, Table 4-2 also shows the

point-based preliminary remediation goals (PRGs) for total PCBs and BaPEq (Tetra Tech, 2013c).

The point-based PRG and point-based RAL are the same for total PCBs (0.676 milligrams per kilogram (mg/kg), whereas the PRG for BaPEq (0.70 mg/kg) is approximately one order of magnitude (i.e., a power of 10) lower than the RAL (6.50 mg/kg). The Middle River Complex (MRC) sediment feasibility study developed PCB RALs to be protective of both human health and ecological receptors at the site. The RALs were developed using United States Environmental Protection Agency (USEPA)-approved methodologies that incorporate the most current scientific knowledge on the effects of these compounds on potential receptors. USEPA considers the ecological RAL, which is the consensus-based freshwater probable-effect concentration developed by MacDonald, et al., (1996, 2000a, b) and adopted by the USEPA Biological Technical Assistance Group (USEPA, 2006a, b), to be protective of benthic invertebrates. Although the human health RAL is lower than the ecological sediment RAL, the human health RAL is an areaweighted concentration that cannot be compared directly to the point-based (i.e., single sample) storm drain system sediment concentrations; therefore the lowest point-based RAL of 0.676 mg/kg developed for ecological receptors was selected to screen the storm drain sediment sample results in this report. Based on the sediment risk assessment, this ecological-based remedial level (0.676 mg/kg) would also not present an unacceptable human health risk if a human receptor were to directly contact the sediments. Additionally, based on the low volume of sediments in the storm drain system, contribution of storm drain sediment to Dark Head Cove at concentrations at or below 0.676 mg/kg would not substantially contribute to the areal extent of PCB in sediment in the receiving waters (i.e., not substantially raise the area-wide PCB concentration).

Table 4-2 and Figure 4-1 show total PCB and BaPEq sampling results that exceed the proposed point-based sediment RALs (0.676 milligrams per kilogram [mg/kg] and 6.50 mg/kg, respectively). Total PCB and BaPEq concentrations in these tables are presented in units of both milligrams per kilogram and micrograms per kilogram (μg/kg), for consistency with previous MRC documents.

#### 4.1 POLYCHLORINATED BIPHENYLS

PCBs were detected in all 32 sediment samples (Tables 4-1 and 4-2), at concentrations ranging from 0.0021 mg/kg (sample SD-170-0–1 in the MRC Outfall 004 system) to 780 mg/kg (sample SD-182-0–3 in the MRC Outfall 005 system). As shown in Table 4-2 and Figure 4-1, total PCB concentrations exceed the proposed sediment RAL (0.676 mg/kg) in 10 sediment samples within five outfall systems (MRC outfall systems 005, 007, 008, and 009, and Martin State Airport [MSA] outfall system IN159). Samples exceeding the proposed RAL (0.676 mg/kg) are summarized in Table 4-3.

Sediment samples with the highest concentrations of total PCBs (40–780 mg/kg) were collected in MRC Block E, in the Outfall 005 and 008 systems in the southern portion of the property. One elevated PCB concentration (SD-171-0–1 at 150 mg/kg) was detected at the northern edge of Block F, but it is directly downstream of catch basins and manholes where several of the highest PCB concentrations are present in Block E (e.g., SD-182-0–3 [780 mg/kg], SD-173-0–3 [220 mg/kg], and SD-172-0–3 [54 mg/kg]. Lower exceedances were detected at concentrations ranging from 0.74–1.71 mg/kg, in samples collected southeast and east of MRC Building C (samples SD-185-0–1 and SD-191-0–1) and northeast of the Hangar 4–6 building at Martin State Airport (MSA) (sample SD-175-0–1).

#### 4.2 BENZO(A)PYRENE EQUIVALENT

PAHs were detected in all 32 sediment samples (Tables 4-1 and 4-2), with BaPEq concentrations ranging from 0.098 mg/kg (sample SD-169-0–1 in MRC Outfall 003 system) to 64.0 mg/kg (sample SD-190-0–1 in MRC Outfall 007 system). As shown in Table 4-2 and Figure 4-1, BaPEq concentrations exceeded the proposed sediment RAL (6.50 mg/kg) in seven sediment samples within five outfall systems: MRC Outfall systems 005, 007, and 008, and MSA Outfall systems 90F001 and IN159.

Sediment samples with BaPEq concentrations greater than the proposed sediment RAL (i.e., BaPEq concentrations ranging from 8.06–64.0 mg/kg) were collected at the following locations (listed from highest to lowest concentration):

• The MRC Outfall 007 system, in the catch basins east of Building C (samples SD-190-0-1 and SD-191-0-1, at 19.1 and 64.0 mg/kg, respectively). These catch basins

contain only approximately one inch of sediment each, and are at the upstream end of the Outfall 007 system. Transport of this sediment to the Dark Head Cove outfall is unlikely, because of the limited quantity of sediment in the catch basins, the long travel distance to the outfall, and the absence of sediment in manholes inspected downstream (i.e., northeast and southeast) of these catch basins.

- MSA Outfall 9OF001, in an open channel along Wilson Point Road (sample SD-177-0-2 at 30.5 mg/kg). This sample was collected from an open channel along Wilson Point Road where transport of sediment is possible to downstream locations; however, the BaPEq concentration of 2.70 mg/kg in downstream sample SD-176-0-3 indicates that PAHs in sediment are not likely reaching Dark Head Cove at concentrations above the RAL (6.50 mg/kg).
- The MRC Outfalls 005 and 008 systems in MRC Block E (samples SD-181-0-1, SD-182-0-3, and SD-183-0-1, range of 14.8-25.5 mg/kg). Sediment at these three locations is unlikely to be transported to Dark Head Cove because of an elevated outflow pipe in manhole MH-9 (location of SD-181-0-1). MH-9 is downstream of the other two sampling locations that had elevated BaPEq in this system (SD-182-0-3 and SD-183-0-1). The higher outflow pipe at MH-9 is usually partially filled with water, and has an area where sediment drops out and is trapped in MH-9 upstream of the outfall. The presence of only approximately one inch of sediment at MH-9 indicates that sediment is trapped in MH-9 (i.e., the accumulated sediment level in MH-9 is below the outflow pipe), and is likely not reaching Dark Head Cove.
- The manhole at the northeastern corner of MSA Hangar 4 (sample SD-175-0–1, at 8.06 mg/kg) contains only approximately one inch of sediment and is near the upstream end of the Outfall IN159 system. Transport of this sediment to the outfall is unlikely due to the limited quantity of sediment in the manhole, the long travel distance to the outfall, and the absence of sediment in manholes inspected downstream of this location. Furthermore, the BaPEq concentration (2.87 mg/kg) in downstream sample SD-174-0–1 indicates that PAHs in sediment are not likely reaching Dark Head Cove at concentrations above the RAL (6.50 mg/kg).

#### 4.3 POSSIBLE SOURCES DISCUSSION

Contaminant concentrations and observed volumes of sediment in the drainage systems indicate that MRC outfall systems 001–004 are not significant continuing sources of PCBs and PAHs (expressed as BaPEq) to Cow Pen Creek sediment, because PCB and BaPEq concentrations in sediment samples of these outfall systems are below the RALs. Additionally, the amount of sediment (one inch or less) noted in these outfall systems is sparse. Outfalls 001–004 primarily receive runoff from building roofs and paved portions of the MRC, where PCBs are not chemicals of concern. This storm-drain-sediment investigation suggests that MRC Outfall 005 may act as a continuing source of PCBs and PAHs (expressed as BaPEq) to Dark Head Cove sediment, because concentrations of these two constituents in sediment samples are substantially

above the RALs (i.e., PCB concentrations up to 780 mg/kg and BaPEq concentrations up to 30.5 mg/kg). Sediment is also available to be transported via the Outfall 005 system.

Several inches of sediment were encountered in two Block E catch basins sampled, and one manhole (sample SD-173 at MH-4) was filled with soil/sediment. The manhole at sampling location SD-173 (MH-4) was full of sediment at the time of the 2011 IRM. Manhole MH-4 and sampling location SD-172-0–3 (in catch basin IL-18) were not cleaned because the pipes are blocked in this part of the Outfall 005 system. Sampling location SD-182-0–3 (manhole MH-7) was cleaned and repaired during the IRM. However, transport of sediment to Dark Head Creek may be limited, since much of this system is still blocked, because discharge from the outfall is low, and because downstream manholes nearest Outfall 005 (e.g., MH-1 and MH-2) contain little to no sediment several years after completion of the IRM.

MSA Outfalls OF9001 and IN159 might possibly act as limited continuing sources of PCBs and PAHs (represented as BaPEq) to sediment of Dark Head Cove, based on upstream sample concentrations of one or both of these constituents above the RALs. However, downstream sediment sampling results (SD-174-0–1 and SD-176-0–1) suggest that concentrations of PCBs and BaPEq above the RALs may not be transported via sediment to Dark Head Cove. Accurate estimates of sediment quantities in the storm drain systems cannot reliably be made from this study.

Except for the MRC Outfall 005 system, samples collected and/or visual inspections conducted for this study indicate only small quantities of sediment in catch basins and manholes. Portions of the Outfall 005 system cleaned during the IRM and sampled for this investigation (samples SD-171-0–1, SD-181-0–1, and SD-182-0–3) contain only limited quantities of sediment (thicknesses of one to three inches), although sediment concentrations exceed the PCB and/or BaPEq RALs at these locations. The conditions of Outfall 005 manholes and catch basins not cleaned during the IRM but sampled as during this investigation (SD-172-0–3, SD-173-0–3, and SD-183-0–1) appear to be the same as those reported during the 2011 IRM. Previous camera studies at the Middle River Complex (Tetra Tech, 2012a, 2012d) demonstrate that most of the sediment in storm drainage systems accumulates at pipe joints (particularly offset joints or mineral encrustations), where it cannot be observed without the aid of a mobile camera system.

Therefore, underestim					and	manholes	likely

Table 4-1

Statistical Summary of Analytes Detected in Storm Drainage System Sediment Samples-2013

Lockheed Martin Middle River Complex and Martin State Airport

Middle River, Maryland

	Frequency		Minimum	Maximum	Minimum	Maximum	Location of	Sample of	Average of	Average of	Standard
Parameter	of	Percent	Non-Detect	Non-Detect	Detected	Detected	Maximum	Maximum	All Results	Positive	Deviation
	Detection		Concentration	Concentration	Concentration	Concentration	Concentration	Concentration	All Results	Results	Deviation
Polychlorinated biphenyls (mg	/kg)										
TOTAL AROCLOR	32/32	100			0.0021	780	SD-182	SD-182-0-3	52.2	52.2	157.9
Polycyclic aromatic hydrocarb	ons (mg/kg)										
BaPEq-POS	32/32	100			0.096	64.0	SD-190	SD-190-0-1	7.02	7.02	13.1
Polychlorinated biphenyls (µg/	kg)										
AROCLOR-1254	18/32	56	0.37 U	2500 U	5.3	410	SD-175,	SD-175-0-1,	149.3	105.3	289.3
7 IKOCEOK-1254			0.57 0	2300 0	5.5	410	SD-184	SD-184-0-2	147.5	103.3	207.3
AROCLOR-1260	32/32	100			2.1	780000	SD-182	SD-182-0-3	52160	52160	157900
TOTAL AROCLOR	32/32	100			2.1	780000	SD-182	SD-182-0-3	52220	52220	157900
Polycyclic aromatic hydrocarb	ons (µg/kg)										
BaPEq-POS	32/32	100			95.75	64050	SD-190	SD-190-0-1	7023	7023	13100
BENZO(A)ANTHRACENE	32/32	100			76 J	49000	SD-190	SD-190-0-1	4952	4952	9738
BENZO(A)PYRENE	32/32	100			69 J	43000	SD-190	SD-190-0-1	4685	4685	8803
BENZO(G,H,I)PERYLENE	32/32	100			81 J	33000	SD-190	SD-190-0-1	3752	3752	6755
CHRYSENE	32/32	100			120 J	55000	SD-190	SD-190-0-1	5804	5804	11110
FLUORANTHENE	32/32	100			180 J	150000	SD-190	SD-190-0-1	14120	14120	30300
PYRENE	32/32	100			160 J	74000	SD-190	SD-190-0-1	9394	9394	17450
BENZO(B)FLUORANTHENE	31/32	97	130 U	130 U	120 J	47000	SD-190	SD-190-0-1	5650	5830	10090
BENZO(K)FLUORANTHENE	31/32	97	170 U	170 U	43 J	19000	SD-190	SD-190-0-1	2161	2227	3987
INDENO(1,2,3-CD)PYRENE	31/32	97	86 U	86 U	66 J	30000	SD-190	SD-190-0-1	3346	3453	6104
PHENANTHRENE	31/32	97	130 U	130 U	47 J	61000	SD-190	SD-190-0-1	7519	7759	14600
ANTHRACENE	30/32	94	19 U	82 U	30 J	18000	SD-190	SD-190-0-1	1597	1701	3473
DIBENZ(A,H)ANTHRACENE	30/32	94	21 U	48 U	26 J	8200	SD-190	SD-190-0-1	917.6	977.6	1667
ACENAPHTHYLENE	28/32	88	22 U	96 U	14	870	SD-190	SD-190-0-1	185.4	208	198.1
FLUORENE	26/32	81	25 U	110 U	13	6600	SD-190	SD-190-0-1	720.1	880.7	1439
ACENAPHTHENE	25/32	78	20 U	80 U	9 J	5600	SD-190	SD-190-0-1	704.2	895.9	1328
NAPHTHALENE	19/32	59	6.8 U	72 U	4.7 J	640	SD-181	SD-181-0-1	112	179.7	169.5
2-METHYLNAPHTHALENE	18/32	56	7.1 U	75 U	2.5 J	530	SD-190	SD-190-0-1	90.9	151.3	136.6
	•		•	-	-	Associated Samp	oles:	•	•		
BaPEq - benzo(a)pyrene equiv	alent					SD-163-0-1	SD-171-0-1	SD-179-0-1	SD-187-0-1		
J - estimated concentration						SD-164-0-1	SD-172-0-3	SD-180-0-1	SD-188-0-1		
μg/kg - micrograms per kilogr	am (i.e., par	ts per billio	1)			SD-165-0-1	SD-173-0-3	SD-181-0-1	SD-189-0-1		
7 711 171						CD 166 0 1	CD 174 0 1	GD 102 0 2	CD 100 0 1		

BaPEq - benzo(a)pyrene equivalent	SD-163-0-1	SD-171-0-1	SD-179-0-1	SD-187-0-1
J - estimated concentration	SD-164-0-1	SD-172-0-3	SD-180-0-1	SD-188-0-1
μg/kg - micrograms per kilogram (i.e., parts per billion)	SD-165-0-1	SD-173-0-3	SD-181-0-1	SD-189-0-1
mg/kg - milligrams per kilogram (i.e., parts per million)	SD-166-0-1	SD-174-0-1	SD-182-0-3	SD-190-0-1
PCBs - polychlorinated biphenyls	SD-167-0-1	SD-175-0-1	SD-183-0-1	SD-191-0-1
POS - only detected polycyclic aromatic hydrocarbons are used for this calculation	SD-168-0-1	SD-176-0-3	SD-184-0-2	SD-192-0-1
U - not detected at the concentration shown left of the letter	SD-169-0-1	SD-177-0-2	SD-185-0-1	SD-193-0-1
not applicable	SD-170-0-1	SD-178-0-1	SD-186-0-1	SD-194-0-1

Table 4-2
Analytes Detected and Screening Criteria for Storm Drainage System Sediment Samples-2013
Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland
Page 1 of 4

LOCATION	Preliminary	Remedial	SD-163	SD-164	SD-165	SD-166	SD-167	SD-168	SD-169	SD-170
SAMPLE ID	Remedial	Action	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-167-0-1	SD-168-0-1	SD-169-0-1	SD-170-0-1
SAMPLE DATE	Goal <sup>(1)</sup>	Level <sup>(1)</sup>	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/20/2013
Polycyclic aromatic hydrocarbons	(mg/kg)									
BaPEq-POS	0.70	6.50	0.775	0.967	1.43	0.233	0.382	0.593	0.096	0.362
Polychlorinated biphenyls (mg/kg)										
TOTAL AROCLOR	0.676	0.676	0.025	0.089	0.099	0.034	0.019	0.184	0.058	0.0021
Polycyclic aromatic hydrocarbons	(μg/kg)									
BaPEq-POS	700	6500	775.31	966.96	1429.4	233.18	382.14	592.83	95.75	362.06
2-METHYLNAPHTHALENE				30 J	25 J					
ACENAPHTHENE			64 J	25 J	47 J			110 J	30 J	
ACENAPHTHYLENE			69 J	150 J	36 J					58 J
ANTHRACENE			97 J	200 J	110 J		44 J	180 J		110 J
BENZO(A)ANTHRACENE			530 J	520	830	130 J	270 Ј	470	76 J	320
BENZO(A)PYRENE			500 J	660	920	120 J	280 J	390 J	69 J	250 J
BENZO(B)FLUORANTHENE			700 J	640	1400		430	250 J	120 J	320 J
BENZO(G,H,I)PERYLENE			480 J	760	890	130 J	340 J	410 J	81 J	170 J
BENZO(K)FLUORANTHENE			260 J	420	590		170 J	530	43 J	78 J
CHRYSENE			710 J	760	1500	180 J	440	530	120 J	280
DIBENZ(A,H)ANTHRACENE			110 J	130 J	200 J	100 J		88 J		32 J
FLUORANTHENE			1400 J	960	3500	180 J	800	1000	180 J	630
FLUORENE			52 J	52 J	55 J			86 J		
INDENO(1,2,3-CD)PYRENE			390 J	560	790		300 J	370 J	66 J	150 J
NAPHTHALENE				34 J	95 J			36 J		
PHENANTHRENE			990 J	370	2800		340 J	710	47 J	300
PYRENE			1100 J	1100	2500	210 J	590	830	160 J	480
Polychlorinated biphenyls (µg/kg)										
AROCLOR-1254	676	676						100	29	
AROCLOR-1260	676	676	25	89	99 J	34	19	84	29	2.1 J
TOTAL AROCLOR	676	676	25	89	99	34	19	184	58	2.1

Table 4-2
Analytes Detected and Screening Criteria for Storm Drainage System Sediment Samples-2013
Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland
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LOCATION	Preliminary	Remedial	SD-171	SD-172	SD-173	SD-174	SD-175	SD-176	SD-177	SD-178
SAMPLE ID	Remedial	Action	SD-171-0-1	SD-172-0-3	SD-173-0-3	SD-174-0-1	SD-175-0-1	SD-176-0-3	SD-177-0-2	SD-178-0-1
SAMPLE DATE	Goal <sup>(1)</sup>	Level <sup>(1)</sup>	12/19/2013	12/19/2013	12/19/2013	12/19/2013	12/19/2013	12/19/2013	12/19/2013	12/19/2013
Polycyclic aromatic hydrocarbons (	mg/kg)									
BaPEq-POS	0.70	6.50	1.02	6.39	2.83	2.87	8.06	2.70	30.5	0.258
Polychlorinated biphenyls (mg/kg)										
TOTAL AROCLOR	0.676	0.676	150	54	220	0.5	1.71	0.124	0.101	0.0173
Polycyclic aromatic hydrocarbons (	μg/kg)									
BaPEq-POS	700	6500	1020.22	6393	2828.3	2870.2	8061.1	2699.3	30526	258.05
2-METHYLNAPHTHALENE			14 J	50 J	55 J	47 J	210 J	30 J	300 J	
ACENAPHTHENE			28 J	550	150 J	230	1500	130 J	2200 J	
ACENAPHTHYLENE			71	160 J	430	140 J	310	200 J	510 J	48 J
ANTHRACENE			160	1100	580	610	2100	390	3300 J	30 J
BENZO(A)ANTHRACENE			620	4600	1700	2200	5700	1600	18000 J	140 J
BENZO(A)PYRENE			680 J	4200 J	1900 J	1900 J	5400 J	1800 J	21000 J	180 J
BENZO(B)FLUORANTHENE			720 J	5300 J	2100 J	2500 J	6000 J	2700 J	27000 J	200 J
BENZO(G,H,I)PERYLENE			630 J	3300 J	1800 J	1300 J	3700 J	1500 J	15000 J	140 J
BENZO(K)FLUORANTHENE			250 J	1800 J	630 J	780 J	2500 J	710 J	10000 J	86 J
CHRYSENE			720	5000	2000	2400	6100	2200	26000 J	190 J
DIBENZ(A,H)ANTHRACENE			150 J	870 J	390 J	360 J	1100 J	320 J	3500 J	32 J
FLUORANTHENE			910	9600	3000	5800	13000	4200	68000 J	260
FLUORENE			35 J	460	160 J	250	1000	160 J	2200 J	
INDENO(1,2,3-CD)PYRENE			530 J	3100 J	1500 J	1300 J	3600 J	1400 J	14000 J	110 J
NAPHTHALENE			24 J	89 J	94 J	160 J	480	51 J	120 J	
PHENANTHRENE			430	5000	1100	3800	9000	1700	39000 J	81 J
PYRENE			1100	7400	3400	4000	10000	3000	45000 J	210
Polychlorinated biphenyls (µg/kg)				•		•	•	•	•	
AROCLOR-1254	676	676				150	410	36	28 J	5.3
AROCLOR-1260	676	676	150000	54000	220000	350	1300	88	73 J	12
TOTAL AROCLOR	676	676	150000	54000	220000	500	1710	124	101	17.3

Table 4-2
Analytes Detected and Screening Criteria for Storm Drainage System Sediment Samples-2013
Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland
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LOCATION	Preliminary	Remedial	SD-179	SD-180	SD-181	SD-182	SD-183	SD-184	SD-185	SD-186
SAMPLE ID	Remedial	Action	SD-179-0-1	SD-180-0-1	SD-181-0-1	SD-182-0-3	SD-183-0-1	SD-184-0-2	SD-185-0-1	SD-186-0-1
SAMPLE DATE	Goal <sup>(1)</sup>	Level <sup>(1)</sup>	12/19/2013	12/19/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013
Polycyclic aromatic hydrocarbons	(mg/kg)									
BaPEq-POS	0.70	6.50	0.289	4.32	25.5	14.8	20.1	5.97	1.86	0.471
Polychlorinated biphenyls (mg/kg)										
TOTAL AROCLOR	0.676	0.676	0.0296	0.116	420	780	40	1.71	0.74	0.098
Polycyclic aromatic hydrocarbons	(/)									
BaPEq-POS	(μg/kg) 700	6500	289.43	4322.7	25508	14776	20148	5971.9	1857.3	471.02
2-METHYLNAPHTHALENE	700	0300	2.5 J	4322.7	23308 440 J	240 J	280	170	21 J	4/1.02
ACENAPHTHENE			18	70 J	3900	1600	3000	920	53 J	
ACENAPHTHYLENE			14	76 J	350 J	230 J	320	560	94 J	50 J
ANTHRACENE			41	390 J	7100	3000	5600	2200	230	96 J
BENZO(A)ANTHRACENE			120	2600	20000	10000	15000	4500	1300 J	280 J
BENZO(A)PYRENE			200 J	2900	17000	9500	13000	3800	1200 J	290 Ј
BENZO(B)FLUORANTHENE			290 J	4300	21000	13000	15000	4700	1500 J	540
BENZO(G,H,I)PERYLENE			150 J	2600	14000	9200	11000	3400	1100	350 Ј
BENZO(K)FLUORANTHENE			120 J	1900	8800	3500	6200	1700	570	150 J
CHRYSENE			230	3700	20000	11000	16000	4900	1600	520
DIBENZ(A,H)ANTHRACENE			32 J	490	3100	2200	3100	930	270	67 J
FLUORANTHENE			410	6200	35000	18000	33000	11000	3600	1100
FLUORENE			13	110 J	4100	1200	2400	1100	87 J	
INDENO(1,2,3-CD)PYRENE			150 J	2200	12000	7300	9700	3000	1000	300 J
NAPHTHALENE			4.7 J		640	450 J	480	340	43 J	
PHENANTHRENE			150	2600	28000	12000	17000	6400	1100 J	400
PYRENE			290	5900	37000	18000	19000	6100	1900	590
Polychlorinated biphenyls (µg/kg)										
AROCLOR-1254	676	676	7.6	21				410	190	38
AROCLOR-1260	676	676	22	95	420000	780000	40000	1300	550	60
TOTAL AROCLOR	676	676	29.6	116	420000	780000	40000	1710	740	98

Table 4-2
Analytes Detected and Screening Criteria for Storm Drainage System Sediment Samples-2013
Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland
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LOCATION	Preliminary	Remedial	SD-187	SD-188	SD-189	SD-190	SD-191	SD-192	SD-193	SD-194
SAMPLE ID	Remedial	Action	SD-187-0-1	SD-188-0-1	SD-189-0-1	SD-190-0-1	SD-191-0-1	SD-192-0-1	SD-193-0-1	SD-194-0-1
SAMPLE DATE	Goal <sup>(1)</sup>	Level <sup>(1)</sup>	12/20/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013
Polycyclic aromatic hydrocarbons (	(mg/kg)									
BaPEq-POS	0.70	6.50	2.17	0.954	4.95	64.0	19.1	0.220	0.251	0.188
Polychlorinated biphenyls (mg/kg)										
TOTAL AROCLOR	0.676	0.676	0.121	0.064	0.093	0.099	0.88	0.033	0.017	0.016
Polycyclic aromatic hydrocarbons (	(μg/kg)									
BaPEq-POS	700	6500	2168.9	954.21	4948.7	64045	19123	219.73	251.07	188.49
2-METHYLNAPHTHALENE			48 J			530	230			
ACENAPHTHENE					130 J	5600	2000	20 J	13 J	9 J
ACENAPHTHYLENE			190 J	120 J	160 J	870	440	89	31 J	48 J
ANTHRACENE			260 J	140 J	590	18000	4200	89	54 J	43 J
BENZO(A)ANTHRACENE			1100	470	3000	49000	13000	130	160	100
BENZO(A)PYRENE			1400	620	3300	43000	13000	150	170	130
BENZO(B)FLUORANTHENE			1800	800	3900	47000	16000	170	200	140
BENZO(G,H,I)PERYLENE			1600	690	2800	33000	9200	120	130	94
BENZO(K)FLUORANTHENE			720	250 J	1500	19000	5600	64 J	87 J	44 J
CHRYSENE			1700	710	3700	55000	17000	190	200	150
DIBENZ(A,H)ANTHRACENE			340 J	150 J	680	8200	2300	29 Ј	33 J	26 J
FLUORANTHENE			2900	1000	8100	150000	67000	360	440	240
FLUORENE			82 J	43 J	180 J	6600	2400	34 J	23 J	15 J
INDENO(1,2,3-CD)PYRENE			1300	540	2600	30000	8500	99	110	79 J
NAPHTHALENE				-	66 J	120 J	87 J			-
PHENANTHRENE			850	260 J	2700	61000	42000	170	160	81
PYRENE			1900	710	4500	74000	49000	240	230	160
Polychlorinated biphenyls (µg/kg)			•		•	•	•	•	•	
AROCLOR-1254	676	676	29	15	11	53	350	13		
AROCLOR-1260	676	676	92	49	82	46	530	20	17	16
TOTAL AROCLOR	676	676	121	64	93	99	880	33	17	16

1) Point-based preliminary remediation goals and remedial action levels are from Table 5-1 of the Middle River Complex sediment feasibility study (Tetra Tech, 2013a).

Shaded cell indicates the concentration exceeds the remedial action level.

BaPEq - benzo(a)pyrene equivalent

J - estimated concentration

μg/kg - micrograms per kilogram (i.e., parts per billion)

mg/kg - milligrams per kilogram (i.e., parts per million)

PCBs - polychlorinated biphenyls

POS - only detected polycyclic aromatic hydrocarbons are used for this calculation

-- not detected

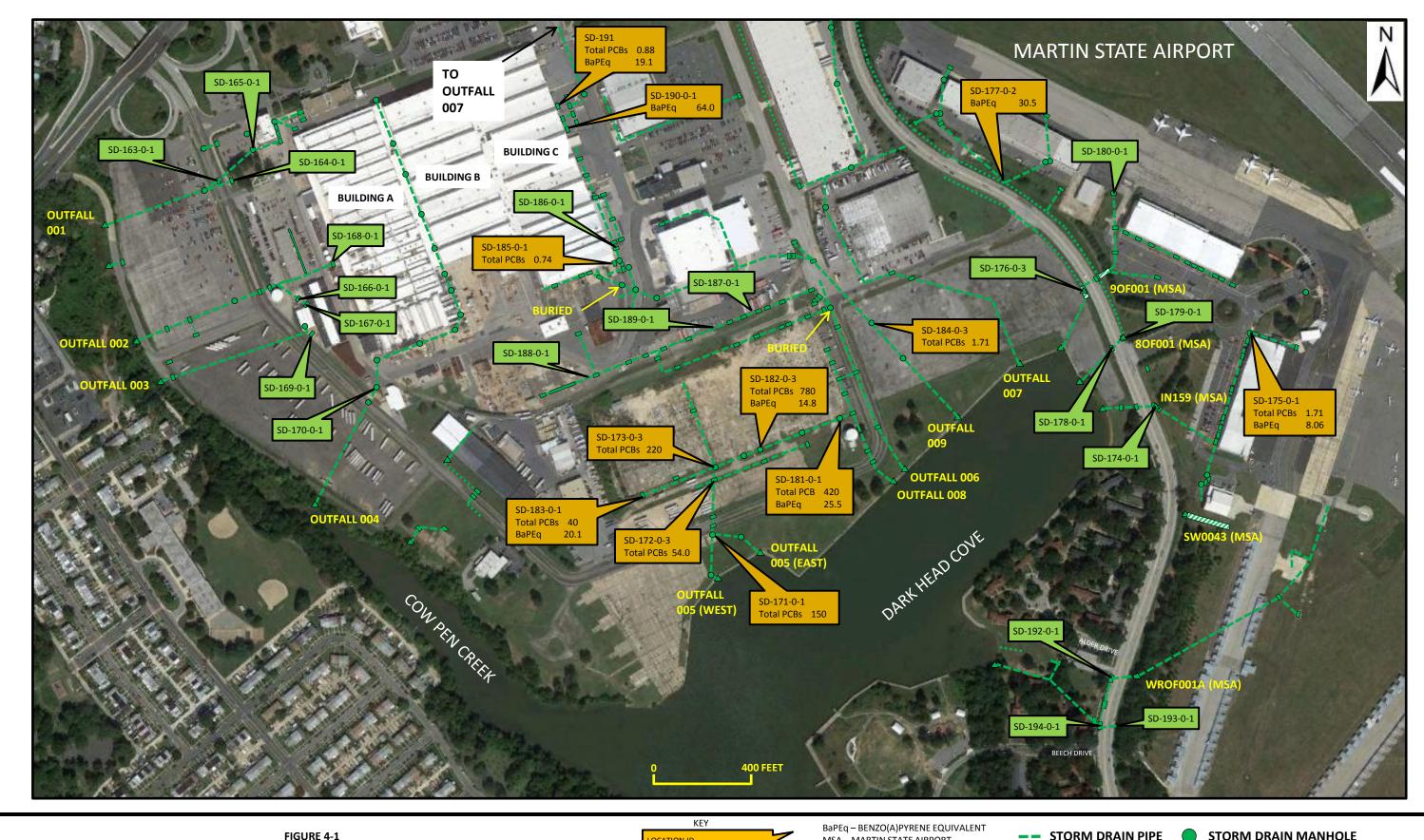
Table 4-3
PCBs in Storm Drainage System Sediment Samples Exceeding the Sediment Remedial Action Level\*—2013
Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland

Sample ID	Total PCB concentration (mg/kg)	Location
SD-182-0-3	780	MRC Outfalls 005/008 systems; along the main line east of the south–central portion of MRC Block E
SD-181-0-1	420	MRC Outfall 008 system; along the main line in the southeastern portion of MRC Block E, near the water tower
SD-173-0-3	220	MRC Outfall 005 system; at the intersection of two main lines in south–central portion of MRC Block E
SD-171-0-1	150	MRC Outfall 005 system; manhole adjacent to roadway near the north side of MRC Block F
SD-172-0-3	54.0	MRC Outfall 005 system; main line in the south-central portion of MRC Block E; sample collected near ground surface in a manhole backfilled with sediment/soil
SD-183-0-1	40	MRC Outfall 005 system; upstream catch basin in the main line in the southwestern portion of MRC Block E
SD-184-0-3	1.71	MRC Outfall 009 system; covered manhole in the main line in the northwestern portion of Lot 6 in MRC Block D; downstream of SD-185-0-1
SD-175-0-1	1.71	MSA Outfall IN159 system; northwestern corner of Hangar 4
SD-191-0-1	0.88	MRC Outfall 007 system; east of Building C in MRC Block I
SD-185-0-1	0.74	MRC Outfall 009 system; southeastern corner of Building C in MRC Block I

\*Remedial action level is the point-based remedial action level of 0.676 for PCBs (Tetra Tech, 2013c).

mg/kg = milligram(s) per kilogram MRC = Middle River Complex MSA = Martin State Airport PCB = polychlorinated biphenyl

SD = sediment



PCBs and BaPEq IN STORM DRAINAGE SYSTEM SEDIMENT SAMPLES EXCEEDING SEDIMENT REMEDIAL ACTION LEVELS -2013 MIDDLE RIVER COMPLEX AND MARTIN STATE AIRPORT, MIDDLE RIVER, MARYLAND

BASED ON TAI 2002 UTILITY PLAN, MARTIN STATE AIRPORT UTILITY COVERAGE, AERIAL PHOTOGRAPHS, TETRA TECH FIELD SURVEYS, AND TETRA TECH SITE INSPECTIONS; ALL UTILITIES BASED ON TAI 2001 UTILITY PLAN. ALL MARTIN STATE AIRPORT COVERAGE NOT VERIFIED IN THE FIELD

OCATION ID CONCENTRATION ANALYTE

GREEN SHADING - NO EXCEEDANCES CONCENTRATIONS ARE IN MILLIGRAMS PER KILOGRAM (MG/KG)

BaPEq - BENZO(A)PYRENE EQUIVALENT MSA – MARTIN STATE AIRPORT PCBs - POLYCHLORINATED BIPHENYLS

**REMEDIAL ACTION LEVELS:** 

BaPEq 6.50 MG/KG PCBs 0.676 MG/KG STORM DRAIN PIPE

OUTFALL

•••• OPEN CHANNEL

STORM DRAIN MANHOLE



STORM WATER INFILTRATION TRENCH

# Section 5 Summary

The following summarizes Lockheed Martin Corporation's (Lockheed Martin's) 2013 storm-drainage-system sediment investigation and findings:

- December 18–20, 2013, Lockheed Martin collected 32 sediment samples from nine Middle River Complex (MRC) storm drainage systems and four Martin State Airport (MSA) storm drainage systems that discharge to Cow Pen Creek or Dark Head Cove.
- Samples were collected to support the remedial design for sediments in Cow Pen Creek and Dark Head Cove. Sediments in these two water bodies contain concentrations of polychlorinated biphenyls (PCBs) and polycyclic aromatic hydrocarbons (PAHs) in excess of proposed point-based (i.e., single sample) remedial action levels (RALs) developed in the feasibility study. The samples were collected to evaluate if sediment in the storm drain systems could potentially act as continuing sources of these two contaminants to sediment in the creek and cove following remediation.
- Most catch basins and manholes inspected lacked sediment or lacked sufficient sediment
  for a sample aliquot. Most appurtenances sampled contained one inch or less of sediment.
  One manhole (sample SD-173-0-3) was backfilled with soil. The field crew could only
  estimate the presence of sediment by using sampling tools and probes in flooded catch
  basins or manholes.
- Sediment samples were chemically analyzed at on off-site laboratory for polychlorinated biphenyls, polycyclic aromatic hydrocarbons, and moisture content.
- Polychlorinated biphenyls and polycyclic aromatic hydrocarbons were detected in all 32 samples collected from the 13 storm drain systems.
- Concentrations of polychlorinated biphenyls exceed the proposed point-based remedial action level of 0.676 milligrams per kilogram (mg/kg) in 10 samples. Polychlorinated biphenyls concentrations above the proposed remedial action level range from 0.74-780 mg/kg.
- The highest concentrations of polychlorinated biphenyls (40–780 mg/kg) were detected in six sediment samples in the Middle River Complex Outfalls 005 and 008 systems in Blocks E and F.
- Concentrations of polycyclic aromatic hydrocarbons (expressed as benzo(a)pyrene equivalents [BaPEq]) in seven samples exceed the proposed point-based benzo(a)pyrene

equivalent remedial action level of 6.50 mg/kg. Concentrations of benzo(a)pyrene equivalent above the proposed remedial action level range from 8.06–64.0 mg/kg, and were detected in samples from the following locations (from highest to lowest concentrations):

- o The MRC Outfall 007 system, in catch basins east of Building C (two samples)
- The MSA Outfall 9OF001 system, in an open channel along Wilson Point Road (one sample)
- o The MRC Outfalls 005 and 008 systems in Block E (three samples)
- The northeastern corner of Martin State Airport Hangar 4 (one sample)
- The investigation indicates that Middle River Complex Outfall systems 001–004 are not significant continuing sources of polychlorinated biphenyls and polycyclic aromatic hydrocarbons to Cow Pen Creek. This conclusion is based on polychlorinated biphenyl and polycyclic aromatic hydrocarbon concentrations below the remedial action levels in these storm-drain-sediment samples, and the sparse amount of sediment (one inch or less) noted in these outfall systems.
- The Middle River Complex Outfall 005 system may be a continuing source of polychlorinated biphenyls and polycyclic aromatic hydrocarbons to the sediment of Dark Head Cove, because concentrations of these constituents are substantially above remedial action levels and sediment is available for transport to Dark Head Cove. However, transport of sediment to Dark Head Creek may be limited, because most of this system is still blocked. Discharge from the outfall is also low, and the downstream manholes nearest Outfall 005 (e.g., MH-1 and MH-2) contain little to no sediment several years after completion of the IRM.
- Martin State Airport Outfalls OF9001 and IN159 may act as limited continuing sources
  of polychlorinated biphenyls and polycyclic aromatic hydrocarbons to sediment of Dark
  Head Cove, because upstream-sample concentrations of one or both of these constituents
  are above the remedial action levels. However, sediment samples collected downstream
  of elevated samples but upstream of the outfalls suggest that sediment containing these
  constituents above the remedial action levels may not be transported to Dark Head Cove.
- Accurate estimates of sediment quantities in the storm drain systems cannot reliably be made from this study. Except for the Middle River Complex Outfall 005 system, sample collection and/or visual inspections conducted for this study indicate only small quantities of sediment in catch basins and manholes. However, previous camera studies at the Middle River Complex (Tetra Tech, Inc. [Tetra Tech], 2010, 2012a, 2012d) demonstrate that sediment typically accumulates at pipe joints (particularly offset joints or mineral encrustations), where it cannot be seen without the aid of a mobile camera system. Therefore, estimates based solely on information from catch basins and manholes likely underestimate the amount of sediment in the storm drainage systems.

Future plans for sediments include additional sediment sampling and chemical analyses of sediments near Outfalls 005 in 2014 to refine the magnitude and extent of more

impacted sediments in support of a remedial design for an Interim Remedial Action to remove elevated concentrations of polychlorinated biphenyls (Note: This future action is now described as the "Outfall 005 Sediment Removal Action"). Complete remediation of Dark Head Cove and Cow Pen Creek where polychlorinated biphenyls, polycyclic aromatic hydrocarbons, and metals exceed proposed cleanup goals is currently being designed and permitted. Future cleanup of Tax Block E soils will include removing polychlorinated biphenyls present in the storm drain system in Block E.

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APPENDIX A—FIELD SAMPLE SHEETS



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nam Project No.: [] Surface Soi [] Subsurface	il	Lockheed Martin Midd 112IC05220 Storm S		Sample I Sample I Sampled C.O.C. N	ocation: SD By: F. k	163-0-1 0-163-0-1 Kolberg	<u> </u>	
[X] Sediment [] Other: [] QA Sample				Type of Sample: [] Low Concentration [] High Concentration				
GRAB SAMPLE DATA	A:							
Date:	12//2/2013	Depth Interval	Color	Descriptio	n (Sand, Silt, C	lay, Moisture, etc.)		
Time: //55			9. 44	SAND.	muist.	Organiz 5		
Method: Bow !	Spuer	0-1	BRUVA			3. <b>3</b>		
Monitor Reading (ppm	THE PERSON NAMED IN		L	5mall	Rucks		$\dashv$	
COMPOSITE SAMPL	E DATA:	KHOHMARKAMERIKAN					000	
Date:	Time	Depth Interval	Color			lay, Moisture, etc.)		
		0"-1"	LT Brown + Black	Gravel	ASAND, 91	raded, Ang	JAN	
Method:				Trace.	EINES	. ,		
				Gravels	uptoty	1		
Monitor Readings				F-C 30	ND	548658		
(Range in ppm):					27.200	202		
N/A				Tr.				
1477								
	_						$\dashv$	
SAMPLE COLLECTION	N INFORMA	TION	Himalogoniaco y parazintación				misorii.	
OAIII EE GOLLEGII	Analysis		Container Requ	irements	Collecte	ed Other		
PAHs, PCBs, Moisture			1 - 8-oz wide mouth gla		(Yes)	No Other	$\neg$	
7 Title, 7 CES, Molecure	Johnson		. O OZ WIGO WIGGIN GI		100			
The Land						=		
OBSERVATIONS / NO				MAP:	wenderen with		mino	
Circle one: MRC r MS								
Outfall Number: OD				L	ot a			
Manhole or CB Numbe	er:			_	•		- 1	
Manhole or Catch Bas		on:						
	Brick/mortar	)						
Dimensions: 35"		1		Ν	NArtin BI	vd.		
Sediment thickness (a Sediment area (approx	pprox):	inches			44443			
Sediment area (approx	V) INDPA	DAT DINZ		~	1			
	Flow	,	_	-	Ð			
	יסיין	•						
	\	/	\		•	7-1		
',	134	. \	)	8			1	
\ \ \	con	cretc \	/				3	
	ROU	uk		ij	ETI CA	te A		
	-	_				10 J		
Circle if Applicable:				Signature(s):			$\neg$	
MS/MSD	Duplicate I	D No.:			ETK			
						part and the second second		



Project Site Nam Project No.: [] Surface So [] Subsurface [X] Sediment [] Other: [] QA Sample	ill e Soil	Lockheed Martin Midd 112IC05220 Storm S		Sample ID No.: Sample Location: Sampled By: C.O.C. No.:  Type of Sample: [] Low Concentration [] High Concentration				
GRAB SAMPLE DAT	A:							
Date:	12//8/2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Moi	sture, etc.)		
Time: 1210			BRUNN		sand, muist	(A)		
Method: Thur (		0.1	BRUWN		•			
Monitor Reading (ppn	n): <b>U.U</b>			dyanic	1			
COMPOSITE SAMPL	.E DATA:							
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Moi	sture, etc.)		
				SAMM 441	It - UF-F SAND	Ludh our		
Method:				in canal	Alak Danah	101		
Metriod.	<b></b>	2) 11 Wester		1	black, organ	gransa		
				10000	DIACK, DIGAN	12		
Monitor Readings				1				
(Range in ppm):				<u> </u>				
N/A		1						
						9		
PAHs, PCBs, Moistur	Analysis e content		Container Req 1 - 8-oz wide mouth o		Collected Yes No	Other		
OBSERVATIONS / N	OTES:			MAP:		and deposit the same		
Circle one MROor M	-			1		HILLIAN HOLDEN		
Outfall Number: 00				Side	malk 1			
Manhole or CB Numb	Table 1			1	. A			
Manhole or Catch Bas	_	n:		1	1 4	Gute A		
Material: Conc.				10. 60		Gute A		
Dimensions: / 2.	×24" V	6" DEEP		1000 100	( )W			
Dimensions: / 2.5 Sediment thickness (a	approx):	inches		Blvd	1			
Sediment area (appro	x): /2" x 2"	וו בצייע		1 4	100 000	<i>14/1</i> 1 1		
	10. 0.0			Martind Blood	1	$\Lambda$		
¥1					G-PASS			
6"	10	+			相			
4.17	10	n· /	1	1	18			
4	8	1	)	1	1			
	10 m			1				
L								
Circle if Applicable:				Signature(s):	11			
MS/MSD	Duplicate IC	No.:		V/		ì		
				1 12/	8	1		



Project Site Nan Project No.:	ne:	Lockheed Martin Midd 112IC05220 Storm S		Sample ID Sample Lo	ocation: SD- 165	-1
[] Surface So [] Subsurface [X] Sediment [] Other: [] QA Sample	e Soil :			C.O.C. No.:  Type of Sample: [] Low Concentration [] High Concentration		
GRAB SAMPLE DAT	`A:					
Date:	12/18/2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Moistu	re, etc.)
Time: 1235 Method: GRAB		0-1"	Brown	SANO: F-	W SAND W/SELT, /U	rsc
Monitor Reading (ppn				WET	poorly graded	
COMPOSITE SAMPL	E DÀTA:					
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Moistu	re, etc.)
Method:				•		
Monitor Readings						
(Range in ppm):	2 2 2 2 2					
N/A						
PAHs, PCBs, Moistur	Analysis e content		Container Requ		Collected (Yes) No	Other
OBSERVATIONS / N				MAP:		
Outfall Number: 60				MAN	TEN BIVA	129
Manhole or CB Numb				The second second		
Manhole or Catch Bas	sin Construction	n:				
	Brick/mortar			, ,	Arti 1	up (1950)
Dimensions: 36 4					my Lot	3 3
Sediment thickness (a Sediment area (appro						
Sediment area (appro		2 8 344	NOSNA WATER	100000	Arking Lot	
	36°	me kn			P' BIL	
	RC	P P-MA	122			
Circle if Applicable:				Signature(s):		
MS/MSD	Duplicate II	O No.:			K	



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nam Project No.: [] Surface So [] Subsurface [X] Sediment [] Other: [] QA Sample	oil e Soil t	Lockheed Martin Midd 112IC05220 Storm S		Sample ID No.: Sample Location: Sampled By: C.O.C. No.:  Type of Sample: [] Low Concentration [] High Concentration
GRAB SAMPLE DAT	_	Danah Intarval	Color	Description (Sand Silt Clay Mainture etc.)
Date: Time:	12/18/2013	Depth Interval		Description (Sand, Silt, Clay, Moisture, etc.)
Method: (-CA)		D"-1"	Black	Black gravel; sand + sitt
Monitor Reading (ppn				graded.
COMPOSITE SAMPL	LE DATA:			
Date:	Time	Depth Interval	Color	Description (Sand, Silt, Clay, Moisture, etc.)
Method:				
Monitor Readings	1 1			
(Range in ppm): N/A				
SAMPLE COLLECTI	ON INFORMAT	TION:		
	Analysis		Container Requ	
PAHs, PCBs, Moistur	e content		1 - 8-oz wide mouth g	glass (Yes) No
_			*	
				C) 50-166
				4
OBSERVATIONS / N	OTES:			MAP:
Circle one MRC or M	SA			1 0/00/1/4
Outfall Number: 💍 🔊	7			* Of-50-166 >
Manhole or CB Numb	er:			
Manhole or Catch Bas	sin Construction	n:		
	Brick/mortar	Contract with the contract of		T SPINTA DROP
		und 3.2 Dec	£	_
Sediment thickness (a		inches and men	t mostly remove	4 4 6 111
Sediment area (appro	)x): 💍	with san	You	-
			4 ±35 4	†
	8 9 s 8 s		Pinet The	tence
Circle if Applicable:			<i>J</i> , -	Signature(s):
MS/MSD	Duplicate ID	No.:		THE



			heed Martin Middle River Complex C05220 Storm Sewer Sediment		Sample ID No.: Sample Location: Sampled By: C.O.C. No.:  Type of Sample: [] Low Concentration [] High Concentration		
	- Vov						
Date:	12/17/2013	Depth Interval	Color		n (Sand, Silt, Clay, Mois	240	
Time: 15:10.		0-1"	40-10	SANO ST	LT Wet wioga,	nied	
Method: GRAB Monitor Reading (pp	m): •//	0-1	Brown	SANO, SILT, WET WOJANUS medgravals			
COMPOSITE SAMP		Carlot III and Carlot III		-	77100		
Date:	Time	Depth Interval	Color	Description	n (Sand, Silt, Clay, Mois	sture, etc.)	
Method:			0 45.155 0 15.155 15.1544				
Monitor Readings		0 Pa	3,233				
(Range in ppm): N/A							
PAHs, PCBs, Moistu	Analysis re content		Container Req 1 - 8-oz wide mouth g		Collected  (Yes No	Other	
OBSERVATIONS / N Circle one MRC or M Outfall Number:  Manhole or CB Numb Manhole or Catch Ba Material: Conc.  Dimensions:  Sediment thickness (i) Sediment area (appro	er: sin Constructio Brick/mortar  "Diam x approx): [1] xx: 28 " x 1"		nte A N N Plastic pu		50-163-0 · ]	Hop inner	
MS/MSD	Duplicate ID	) No.:		Signature(s):	214		



Project Site Name Project No.: [] Surface Soil [] Subsurface S [X] Sediment [] Other: [] QA Sample	Soil	Lockheed Martin Midd			SD-168 By: F. Kolberg	
GRAB SAMPLE DATA:	A SECURE AND A SECURE ASSESSMENT AND A SECURE ASSESSMENT ASSESSMEN					
	2/18/2013	Depth Interval	Color		(Sand, Silt, Clay, Mo	
Time: 154		All . 19	Blacks	5-M STAVE	el w F-B ca	wa mixtur
Method: Grant Monitor Reading (ppm):		0"-1"	DK Brown	wet/		0
COMPOSITE SAMPLE	the state of the s					mond Canania mandoni
		Danah Intanial	0.1	T 5	(O1 O'll Ol M-	
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mo	oisture, etc.)
Method:	- 1	জ্ঞান				
			AUGUSAS			
Monitor Readings		300			100,2772	
			0111700	-		
(Range in ppm):		<del>4.7</del>	198800		10.500	The same of the sa
N/A				l .		A STATE OF THE STA
		2.00		COLUMN TO SEE	18000	
SAMPLE COLLECTION	N INFORMAT	ION:		Weight Mark Control		llari unidullibrii
	\nalysis		Container Req	uirements	Collected	Other
PAHs, PCBs, Moisture	content		1 - 8-oz wide mouth g	lass	Yes No	
4 4 42						
E 1000						
OBSERVATIONS / NOT				MAP:		
Circle one MRO or MSA					2R Tracks	27
Outfall Number: DD				( AST )	- 744	_
Manhole or CB Number:					Hr r	
Manhole or Catch Basin					H	
	rick/mortar				+	
Dimensions: אַייֹגָג		7" Deys		( AST )	- RAMP	Bldg A
Sediment thickness (app		inches				
Sediment area (approx):		"X] "		1	++	
<u>                                      </u>	IN .				H = 1	
	I A		72.		44 1.	ė.
: Est	/\	/·	1 6 0		H ' I	
12" -	14	<b>;</b>	8 6		M = 1	520
out	I N	<b> </b>	• 1		Π -	7
	1. 1.		:/	l	TT	1
	`		<b>/</b>			
No	flow-1	2" Standing "	nate/			
Circle if Applicable:	REALINE DE			Signature(s):		
MS/MSD	Duplicate ID	No.:		1	,	
				14	V	İ
				- / //		
				U		



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Name: Project No.:  [] Surface Soil [] Subsurface Soil [X] Sediment [] Other: [] QA Sample Type:				ocation: SD- log By: F. Kolberg	<u> </u>	
Date:	12/ <b>/\$</b> /2013	Depth Interval	Color	Description	(Sand Silt Clay Mai	oturo eta \
Time: Method:	1610 Glab	6'-1"	Black	Gravel,	I (Sand, Silt, Clay, Moi F-C sand +5ILF IET	Mixtur
Monitor Reading (ppn COMPOSITE SAMPI	n): VA	arregionista de la constitución de			/ p. = (	
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Moi	sture, etc.)
Method:				4	2000 Marin 1990	
Monitor Readings (Range in ppm): N/A						
PAHs, PCBs, Moistur	Analysis		Container Req 1 - 8-oz wide mouth g		Collected (Yes) No	Other
Manhole or CB Numb Manhole or Catch Bas Material: Conc. Dimensions: 23 Sediment thickness (a Sediment area (appro	SA  O3 er: sin Constructio Brick/mortan  × 30" × 2 approx): 2 x): 23" × 3	inches inches  8" d standi	ng water	Signature(s):	Floor diam - Inter-	torage
MS/MSD	Duplicate II	O No.:		[]	1/	



Project Site Nam Project No.:  [] Surface Soi [] Subsurface [X] Sediment [] Other: [] QA Sample	il Soil	Lockheed Martin Midd 112IC05220 Storm S			SD-17D By: F. Kolberg	<i>b-</i> 1
GRAB SAMPLE DATA	Δ.	TED BUTTANDO DE STATEMA DE SE				
Date:	1200/2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture. etc.)
Time: Method:	1515 Cab	0"-1"			, + small stavel wet, loose	
Monitor Reading (ppm COMPOSITE SAMPL	AND DESCRIPTION OF THE OWNER, THE	analan da marena da marena		J	W.1, 1003C	and the second
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
Method:					× 200	
Monitor Readings (Range in ppm): N/A						
PAHs, PCBs, Moisture	Analysis	TION:	Container Requ 1 - 8-oz wide mouth g		Collected  Yes No	Other
Manhole or CB Number Manhole or Catch Bas Material: Conc. Dimensions: 26 Sediment thickness (a Sediment area (approx	SA OOH er: in Constructio Brick/mortar pprox): (): 30 10	inches  12.9 Deep inches	4	MAP:  Manhole 9  SEDSMENT)  Signature(s):	Manhole Sanitary	Conservation Block
MS/MSD	Duplicate II	D No.:		P	Ŋ	



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nat Project No.: [] Surface So [] Subsurfac [X] Sedimen [] Other: [] QA Sampl	oil e Soil t	Lockheed Martin Midd			SD-111 By: F. Kolberg	D-
GRAB SAMPLE DA	TA:					
Date:	12/19/2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture, etc.)
	245	0-1"	Grange-brown	धारी + ८	lmy, loose, wet	
Monitor Reading (ppi	THE RESERVE AND ADDRESS OF THE PARTY OF THE					
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture, etc.)
Method:						
Monitor Readings (Range in ppm): N/A				-		
SAMPLE COLLECT		TION:				
PAHs, PCBs, Moistu	Analysis		Container Requirement 1 - 8-oz wide mouth gla		Collected (Yes) No	Other
TAI 13, TODS, MICISTA	ie content		1 - 0-02 wide modili gia	.33	(Tes) NO	
10 11 11 12 11 12 11 11 11 11 11 11 11 11						
OBSERVATIONS / N				MAP:		Kansan Jahre
Circle one: MRd or M					W	
	05					
Manhole or CB Numb						
Manhole or Catch Ba		n:				
Material: Conc.		0 = 1		C	hesapeake park	مدماع
Dimensions: 29		× 9.51 deep				
Sediment thickness ( Sediment area (appro		inches			Ø 50-171	
Sediment area (appro	JAJ. 10-76	ch	20 10 10 1	- X X		X
		Chesapeal	30."RCP		lark Head Co	v
Circle if Applicable:	BESTERNING BERTH			Signature(s):		
MS/MSD	Duplicate II	O No.:		FIS	•	



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nan Project No.: [] Surface So [] Subsurface [X] Sediment [] Other: [] QA Sample	il Soil	Lockheed Martin Midd 112IC05220 Storm S	***************************************		By: SD-172 F. Kolberg	-3
GRAB SAMPLE DAT	AND DESCRIPTION OF THE PERSON					
Date:	12/ <b>  9</b> /2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture, etc.)
Method: G1		0-3"	Okgney	F-VF SAN	NO W/SELT +SO.	me. Clay
Monitor Reading (ppm COMPOSITE SAMPL	the first warmen and the same of the same	noeseagriniena/Distriction	anonamana wa Waliniya			
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture. etc.)
					***************************************	, , , , , , , , , , , , , , , , , , , ,
Method:						
Monitor Readings						
(Range in ppm):						
N/A		9	<del></del>			-1
, , , ,						2 W 15
SAMPLE COLLECTION	ON INFORMA	TION:				
***************************************	Analysis		Container Requ	uirements	Collected	Other
PAHs, PCBs, Moisture	content	2.02	1 - 8-oz wide mouth gl	lass	(Yes) No	
		- 100 E				
					8 	
				10	50	
OBSERVATIONS / NO	OTES:			MAP:		
Circle one MRC or MS						
Outfall Number: 00				See	E MAP. (IR)	(sama
Manhole or CB Numbe	er: IL-18				•	
Manhole or Catch Bas		n;		TL-18	conside	
Material: Conc.					CONCLOSE	
Dimensions: 27"	xyo"					_
Sediment thickness (a	pprox): 3	inches			1	
Sediment area (approx	O. AT. XAL	6y 3		<i>-</i>	IIII IL-1	4
	$\neg$		_	l (*	1.	•
				<b>\</b>	<b>*</b>	1
		/	\		GRAU	
		(	)			
					Conerete	
Inlant stan		t visible; No fl	<u> </u>			
Circle if Applicable:	- 11.2 MP	N FI	on are stars	Signature(s):		
MS/MSD	Duplicate ID	No.:		FAR	i	
	<u> </u>			(/ 4)		



Project Site Na Project No.: [] Surface Sites S	oil e Soil it	Lockheed Martin Midd 112IC05220 Storm S			Ocation: SD-173 By: F. Kolberg O.:	
				5		
GRAB SAMPLE DA				engeen		
Date:	12/ <b> 9</b> /2013	Depth Interval	Color		(Sand, Silt, Clay, Moisture, e	tc.)
Method: G-	II30	۵"-3"	Brown	F-MSANA Dourly 9	with stulctoy raded	
Monitor Reading (pp		mon. tell alle trender on an		1		
COMPOSITE SAMP			Maria Maria Maria Palata Maria	ND 6 SEATTH SEEDING		
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Moisture, e	tc.)
300						
Method:					200 100	
Monitor Readings						
(Range in ppm):		0		1		
	-					-
N/A						
	<b>——</b>					
					WALLS	
SAMPLE COLLECT		ΠΟN:				
	Analysis		Container Rec		Collected Oth	er
PAHs, PCBs, Moistu	re content		1 - 8-oz wide mouth	glass	(Yes) No	
2000 - 100 -						
		10-10-10-10-10-10-10-10-10-10-10-10-10-1	77. 97.	201-25-12-2		
W.						
OBSERVATIONS / N	IOTES:		s Managaritikan di managari	IMAP:		
Circle one MRC r M						
	05				manhole 50-1	73
Manhole or CB Numb		1		-		
Manhole or Catch Ba				-	J.K.	
Material: Conc.		·····		1	\ /	
Dimensions:		7' (TOP \$ SE	DIMENT)	1	1 5	
Sediment thickness (		inches unkna	wa - Mankele		/ //	
Sediment area (appro	The second secon	METER	un - Mahok 13 fril 06 sedina	]		
most manh	des are 7 <sup>th</sup>	8' deep	o Esedian	<b>A</b>		
		'		<0.04	$\rightarrow$	
				50-174_	→( )~#-¥	
		1	. \		$\gamma$	
		\	<i>'</i>	-5	1	
					IL-18	
	I, & a	HANK-BI DUNTER		1	الماسليل	
NO STRUCTUR	rs (IN/or	standed suster at) vuible -fu	1 of sediment			
Circle if Applicable:				Signature(s):		
MS/MSD	Duplicate ID	No.:		1	11/	l
100000000				1 77	rN	



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Name: Project No.:  [] Surface Soil [] Subsurface Soil [X] Sediment [] Other: [] QA Sample Type:		Lockheed Martin Middle River Complex 112IC05220 Storm Sewer Sediment			Ocation: SD-174 By: F. Kolberg	0-1
GRAB SAMPLE DAT	Δ-					
Date:	12/ <b> 9</b> /2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture. etc.)
Time: 13 W	12/1/12010				d with times	
Method: CHAT	}	0"-1"	Brown to			
Monitor Reading (ppm		Drang-bown	Brown	to orange brow	<b>4</b> 20	
COMPOSITE SAMPL	THE RESIDENCE AND ADDRESS OF THE PERSON NAMED IN					
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
Method:						
Monitor Readings				<u> </u>	A STATE OF THE STA	
(Range in ppm): N/A				W		
SAMPLE COLLECTION INFORMATION:  Analysis  PAHs, PCBs, Moisture content			Container Requirements  1 - 8-oz wide mouth glass  Yes		Yes No	Other
Dimensions: 2 9 Sediment thickness (a Sediment area (appro	SA) V-15 9 er: outfill sin Construction Brick/mortar  I' Diagram approx): /" x): 10 90	inches  B  A  B  A  A  A  A  A  A  A  A  A  A	outall  ox 3.8' Dep	\ \ \ \	13-174 154 154	Parker 4-6 DIX ROAD
MS/MSD	Duplicate II	D No.:		TH		3



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nar Project No.: [] Surface So [] Subsurface	- oil e Soil	Lockheed Martin Mido 112IC05220 Storm S		Sample IE Sample Le Sampled I C.O.C. No	By: SD-195 F. Kolberg	-0-1
[X] Sedimen	ι			Type of Sa	ample: concentration	
[] QA Sample	e Type:				Concentration	14
GRAB SAMPLE DAT						
Date:	12/19/2013	Depth Interval	Color		(Sand, Silt, Clay, Mois	
Time:	14:30	0"-1"	P	M-A SAI	UD, STAT, LOUSE,	, wet
Method:	61ab	0 =1	Brown	2 /	ND, 952T, louse, Maled.	
Monitor Reading (ppr COMPOSITE SAMP	THE RESERVE OF THE PARTY OF THE			Patry	ranea.	
	T		I	THE THE PARTY OF THE PARTY		
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
	1					
Method:					-	
				J		
Monitor Readings		1				7. 1. 1.
(Range in ppm):				All No		
N/A		5.5				
	À			İ		
					, . uxur	
SAMPLE COLLECTI	ON INFORMAT	ION:				
	Analysis		Container Req	uirements	Collected	Other
PAHs, PCBs, Moistui		25.00 25.0000000000000000000000000000000	1 - 8-oz wide mouth g		(Yes) No	
2.00			i iomiomosia ii			
	***					
INTERNAL CL		2				
OBSERVATIONS / N	THE RESERVE AND ADDRESS OF THE PERSON NAMED IN COLUMN TWO			MAP:		
Circle one: MRC of M						
Outfall Number:	IN-15	9				
Manhole or CB Numb				DM C	chel & Henli	AF T
Manhole or Catch Ba		l:		l MA		1
Material: Conc. Dimensions: 26	Brick/mortar	4.75' Dee		· · ·	phole 50-17	75 /\
Sediment thickness (a		inches		i i	DATE IN	37
Sediment area (appro		Inches		1 '	V/	ן ייי ר
	, , , , ,		18" 1	1 4	of 1/11-4/	/
		1	187	l	~	
*		/ •	1 7.	1	4	
	i	1	/ W		/ /HS /	. 1
		1			// /HA	yar 4-b
				TO JN159	N. P. BI	1.
<u> </u>			S RAIN		2/6/	,.,
D		18, 1	Y RAISE			
Circle if Applicable:		Hardel Dubariza Kular		Signature(s):		
MS/MSD	Duplicate ID	No.:			V/	
				1 1/	<i>1</i> 0	



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Name: Project No.:  [] Surface Soil [] Subsurface Soil [X] Sediment [] Other: [] QA Sample Type:		Lockheed Martin Middle River Complex 112IC05220 Storm Sewer Sediment		Sample ID No.: Sample Location: Sampled By: C.O.C. No.:  Type of Sample: [] Low Concentration [] High Concentration	3
GRAB SAMPLE DATA	Δ:				III I I I I I I I I I I I I I I I I I
Date:	12/ 19/2013	Depth Interval	Color	Description (Sand, Silt, Clay, Moisture,	etc.)
	:05		Dk Brown	VF-F SAND + SITT BOOK AC	المل
	RAR	0"_3"	and Die gray	VF-F SAND + SIET, poorly go Dk Brown + Dk Grey love	
Monitor Reading (ppm		2.0. 0/2 1/4 )	Dk Brown + Dk Grey love	Swet	
COMPOSITE SAMPL					
Date:	Time	Depth Interval	Color	Description (Sand, Silt, Clay, Moisture,	etc.)
					,
Nactional.			<u> </u>		
Method:				<u> </u>	
Monitor Readings		Total Borne and Ade	10 11 11 15 15 15 15 15 15 15 15 15 15 15		
(Range in ppm):		- A	1	# Spa	
N/A				- 24 - 24 - 24 - 24 - 24 - 24 - 24 - 24	
SAMPLE COLLECTION	ON INFORMA	TION:			
	Analysis		Container Requ	irements Collected C	ther
PAHs, PCBs, Moisture			1 - 8-oz wide mouth gl		
		9			
- 000 MANUAL - 1 - 100 MANUAL -	5,000 33				
OBSERVATIONS / NO	OTES:			MAP:	CONTRACT S
Circle one: MRCorMS					
Outfall Number: 9	OFOOI	(Downstream)		changel Fence ( E.)	
Manhole or CB Number	er:			Changel Fence [ E ]	x12
Manhole or Catch Bas	in Construction	n:		Mec 11/5/ HAM	<b>*</b>
	Brick/mortar			1 100 1	/
	x 4.2	Deep Box 1	6.5 × 3.9'		u l
Sediment thickness (a	pprox): 3"	inches			Ά Ι
Sediment area (approx	(): 6.5° ×	3.9' x 3"			HAMAT
			_ / / /	Box	<b>1</b>
5			3,4	conerate box	.\
			1/29 N	6.5' Long 3.9' wid	4
		1	4 14	4.5 2014	\
		29"	/ / /	6.5' Long 3.9' wid manble diameter 19"	\
	]	•			V
<u> </u>		Anna de caba d	1 441 241	Depth 4.2'	
A	RFU	correspond	meiri ri		
Circle if Applicable:				Signature(s):	
MS/MSD	Duplicate II	D No.:		1 12 10	
				74	l



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project No.: 112IC0522  [] Surface Soil [] Subsurface Soil [X] Sediment [] Other: [] QA Sample Type:  GRAB SAMPLE DATA:  Date: 12/19/2013 Depth  Time: 510  Method: 6/46  Monitor Reading (ppm): NA		Lockheed Martin Midde 112IC05220 Storm Sto	Sewer Sediment Sample C.O.C.  Type o [] Lov [] Hig		By: SD- 177  F. Kolberg	sture, etc.)
COMPOSITE SAMP	LE DATA:					
Date:	Time	Depth Interval	Color	Description (Sand, Silt, Clay, Moisture		sture, etc.)
Method:						
Monitor Readings (Range in ppm): N/A						
PAHs, PCBs, Moistu	Analysis	ATION:	Container Req 1 - 8-oz wide mouth g		Collected Yes No	Other
OBSERVATIONS / N Circle one: MRC of N Outfall Number: Manhole or CB Numb Manhole or Catch Ba Material: Conc. Dimensions: Sediment thickness (a Sediment area (appro	isa) per: sin Construction Brick/mortar  RCP approx): 2-3  ix ope  dit  HA.			Open dittid	Dark Head Cove	4->
MS/MSD	Duplicate II	O No.:		FA	L.	



Project Site Name: Project No.:  [] Surface Soil [] Subsurface Soil [X] Sediment [] Other: [] QA Sample Type	Lockheed Martin Mid 112IC05220 Storm		Sampled E C.O.C. No Type of Sa [] Low Co	SD- 179 By: F. Kolberg	0-1
GRAB SAMPLE DATA:					
Date: 12/1	7/2013 Depth Interval	Color		(Sand, Silt, Clay, Mois	
Time: (600		N. con	M-C 541	Id w/ K/M 4N.	ا ده افد
Method: GRAB	<u></u> "-1"	Dugray	M-c sand w/ F/M sand and SIET; little fines, lowe, ansvian		
Monitor Reading (ppm):			2151 1141	e tines, lowe,	aviaivi
COMPOSITE SAMPLE DAT	ΓA:				
Date: T	ime Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
Method:					
Monitor Readings			-	345 F-1/400	
Monitor Readings		124	<b>†</b>		
(Range in ppm):		2000	-		
N/A		7 - 7 - 7 - 7 - 7 - 7 - 7 - 7 - 7 - 7 -	ļ	1.300	
	488				
SAMPLE COLLECTION INFORMATION:  Analysis  PAHs, PCBs, Moisture content		100		Collected (Yes) No	Other
OBSERVATIONS / NOTES:			MAP:		
Circle one: MRC of MSA				· · · · · · · · · · · · · · · · · · ·	
Outfall Number: Down 5	tream she 804001		Coste	n basin   Ditch	
Manhole or CB Number:			cute	1614	
Manhole or Catch Basin Cor	nstruction:		1, 150-	178 14 00	79
Material: Conc. Brick/	mortar		1 1/1 1 7	178	overland
Dimensions: 5.6' x 2.7' x 4.3' deep			1 / 6	上即一	- शुक्त
	Sediment thickness (approx): inches				
Sediment area (approx): 30%			MAC 4	pA.S	A
		_	6	Ë	strance
← ← 24" RCP			/ <sub>//</sub> / u	point Rund	-
12.11	0.51 Standing	water	'/	'ROAL \	#
Circle if Applicable:			Signature(s):		
	licate ID No.:		FAX	/	



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nam Project No.:		Lockheed Martin Middle River Complex  112IC05220 Storm Sewer Sediment		Sample ID No.: SD-179-0-1 Sample Location: SD-179 Sampled By: F. Kolberg		
[] Surface Soi [] Subsurface [X] Sediment [] Other: [] QA Sample	Soil					
GRAB SAMPLE DATA	A:	***				
Date:	12/19/2013	Depth Interval	Color		(Sand, Silt, Clay, Moi	
	645		Brown to	SAND + 5.	ET: F-M SAND Me SMAIL Gravels	w/six
Method: G A	AB	0"-1"	orante-planu	and so	me small gravels	, louse, we
Monitor Reading (ppm	):			ALIMITAN DATE OF THE PARTY OF T		Hold in the Management of
COMPOSITE SAMPL						
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Moi	sture, etc.)
Method:						
Monitor Readings		1 180				
(Range in ppm):						
N/A					(4)	
IN/A					XXX (IIII	
			***			
SAMPLE COLLECTION	N INFORMA	TION:		is e la company de la comp		
OAIIII EE GOEEEGTI	Analysis		Container Requ	irements	Collected	Other
PAHs, PCBs, Moisture			1 - 8-oz wide mouth gl		(Yes) No	
		9 (1.17)		giant to the same of the same		
(4133.)		<b>4</b> 1	S- 141,000,1134,0			1
				I Production		
1857E1	10.					
OBSERVATIONS / NO				MAP:		
Circle one: MRC o MS				600	Hardwall 24'	rcp
Outfall Number:	goro	01		NO	1) DHen	
Manhole or CB Number				0.00	LIKE COLUM	
Manhole or Catch Bas		on:		غا ا	Ton Header	made da
	Brick/mortar			A47-62	43.72 . Mines	02
Dimensions: ayı RCP Sediment thickness (approx): J/A inches					Links Under	11 to sample
Sediment area (approx	c): JA		men channel		MSA	
Sediment area (approx): NIA collected from open channel			<b>y 200 200 200</b>	MSA Entranca		
					ENIN	
			WISON	1		
		\	/ /	point, Rosal	\	
			ROAD	1		
		. (	-/11			
sample from open chancel				Clamature (1)		
Circle if Applicable:		D. A.I.		Signature(s):	W.25	1
MS/MSD	Duplicate II	D No.:		F/	U /	
				1	1/	



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nam Project No.: [] Surface Soi [] Subsurface [X] Sediment [] Other: [] QA Sample	il Soil Type:	Lockheed Martin Midd 112IC05220 Storm S		Sample Sample C.O.C.  Type of Low		0-
GRAB SAMPLE DATA	The second second	T Dowth Internal	Color	Dogovina	ion (Sand, Silt, Clay, Moi	oturo eta \
Date: Time: 16	12/ <b>\9</b> /2013 <b>\5</b>	Depth Interval			17 F-000 - 000-00	37.39
Method: Gr	-	0"-1"	Brown to	M-C	sand, gradel, u	n et, 100se
Monitor Reading (ppm	2/477	10-1	DK. Brown			
COMPOSITE SAMPL	NAME OF TAXABLE PARTY.					
Date:	Time	Depth Interval	Color	Descript	ion (Sand, Silt, Clay, Moi	sture, etc.)
Method:			•		1.00	
Monitor Readings						-2/37 12 - 2/97
(Range in ppm): N/A						
SAMPLE COLLECTION INFORMATION:  Analysis  PAHs, PCBs, Moisture content  1 - 8-oz wide mouth g				Collected No	Other	
Manhole or CB Number Manhole or Catch Bas	OFUOI er: in Construction Brick/mortar  11 211	y 3.61 deep		MAP:	MSA HANGAN 1 PArking lot MB-18D	*
Circle if Applicable:				Signature(s		
MS/MSD	Duplicate I	D No.:		f	W	9



#### **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nan Project No.:		Lockheed Martin Midd		Sample ID Sample Lo	ocation: SD-181  By: F. Kolberg	D-1
[] Surface So [] Subsurface [X] Sediment	Soil			C.O.C. No	ample:	
[] Other: [] QA Sample	Type:				oncentration Concentration	
				_ 1,9 4		
GRAB SAMPLE DAT					(0 1 0); O1 41 :	
Date: Time:	12/20/2013 (D:D)	Depth Interval	Color		(Sand, Silt, Clay, Mois	
Method:	KR	0"-1"	Brown	F-c sand, provily graded, 1985, wet, little fines		
Monitor Reading (ppm	the state of the s	,				
COMPOSITE SAMPL	THE PERSON NAMED IN COLUMN 1			AND RESERVED LINES		
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture, etc.)
Method:		7,-1,-33,397.		CONTRACT OF THE PARTY OF THE PA		
Monitor Readings						W-112
(Range in ppm):						
CAMPLE COLLECTI		TON.				
SAMPLE COLLECTION	Analysis	TION:	Container Requ	uiremente	Collected	Other
PAHs, PCBs, Moistur	-		1 - 8-oz wide mouth g		(Yes) No	Other
711107 . 020, moletar	o oomon		1 0 02 mac meaning		100	
OBSERVATIONS / N				MAP:		
Circle one: MRC or M						
Outfali Number:	28 /003				700000	
Manhole or CB Number				1 (	1 MH-9	
Manhole or Catch Bas Material: Conc. (		n:		{	50-181	
Dimensions: 24		x 4.5' Decu		· •		
Sediment thickness (a	A 400	inches				
Sediment area (appro				l		7
	1 212	4	8" pipe	1	Asphalt	
			o Pipo		A	ا ۱
		/ 4	1		(//////////////	, ,
	11	6.	→\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	1	(////) F.	e water
	15" RC	n —	Y RCP N		SUPI	ely
		**************************************	/		TA	ST e water aly alk
	^	1.8 & Standy		l		
		1.8 &Standu	i water			
Circle if Applicable:				Signature(s):		
MS/MSD	Duplicate IC					



# Tetra Tech, Inc. SOIL & SEDIMENT SAMPLE LOG SHEET

Project Site Nam Project No.: [] Surface So [] Subsurface [X] Sediment [] Other: [] QA Sample	il Soil	Lockheed Martin Midd 112IC05220 Storm S			ocation: SD-182 By: F. Kolberg	9-3
GRAB SAMPLE DAT			r a.	T 5	(0l 0'll 0  M-i-	
Date:	12/2013	Depth Interval	Color		(Sand, Silt, Clay, Mois	ture, etc.)
Time:	10:20	0"-3"	Black	STLA + CIA)	, louse, wet	
Method: GP Monitor Reading (ppm		0 -9	TO DK gray			
COMPOSITE SAMPL	MARIE STREET, SQUARE		alemorasionismo medicinasion			Barbara a a a a a a a a a a a a a a a a a
					40 1 011 01 11	STITUTE OF STATES
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture, etc.)
<del></del>						
Method:						
Monitor Readings					0-0	
(Range in ppm):					54.5	
N/A	1		****			
IN/A						
SAMPLE COLLECTION INFORMATION:  Analysis  PAHs, PCBs, Moisture content			Container Req 1 - 8-oz wide mouth g		Collected (Yes) No	Other
	Stores to the control of the	VV.1910-14-15-15-15-15-15-15-15-15-15-15-15-15-15-				
OBSERVATIONS / NO				MAP:		
Circle one: MR) or M	SA					
Outfall Number: 🧿	05				-Block E	
Manhole or CB Number	er: <b>///.ii -7</b>			south	un area	
Manhole or Catch Bas	sin Construction	on:		400	TRM AS-BUI	LT
Material: Conc.		V 1 51 A		1 300	A MANUTUL	
Dimensions: 20	)" NIAA 2 ii	x 37' Deep	0,8 water		91 ACC SNO	N N
Sediment thickness (a Sediment area (appro	ipprox):	t antique ble de	Luc O anda I	•		
Sediment area (appro.	A). Jepine,	ANTOPLE	ant planta	1		
		·				
				1		
İ		4	ااس کھ		(9)	Į.
	L	6"	13	l		
		Rep (	Rep			
	DI	w structures	والمراء ووالمو			
Circle if Applicable:	,	W OTTO MOTO	MAI AIRIA	Signature(s):		
MS/MSD	Duplicate II	D No.:		1	1/1	
	Sapilodio II			1 11	lγ	



Project Site Nam Project No.:	e:	Lockheed Martin Middle River Complex 112IC05220 Storm Sewer Sediment		Sample ID No.: SD-183-0-1 SD-183		
[] Surface Soil				Sampled I	By: F. Kolberg	
[] Subsurface [X] Sediment [] Other: [] QA Sample					ample: oncentration Concentration	
GRAB SAMPLE DATA		ARTHUR MILITER OF THE STREET			TELEVISION NAMED OF THE OWNER, THE OWNER, THE OWNER, THE OWNER, THE OWNER, THE OWNER, THE OWNER, THE OWNER, THE	
	12/20/2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Moi	sture. etc.)
Time: (D:				1	Chay, some san	62
Method: G-R		0'-1'	Black	wet	OM // 3041- 342	n, 10-2-5
Monitor Reading (ppm)	The state of the last of the l					
COMPOSITE SAMPLE				MULICAL DISCONDING		
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Moi	sture, etc.)
Method:					***	
Monitor Readings						
(Range in ppm): N/A					ū	
SAMPLE COLLECTION INFORMATION:  Analysis  PAHs, PCBs, Moisture content			Container Requirements 1 - 8-oz wide mouth g		Collected Yes No	Other
OBSERVATIONS / NO	TEC.			MAP:		
Circle one: MRC or MS				IWAP.		
Outfall Number:	COLU					/
Manhole or CB Numbe Manhole or Catch Basi	に エムーパ			MRC	Bluck E south Area LM AS-BUILT	resid
Material: Conc.		ort.			Area	
Dimensions: 29'x	20" x 4			See II	IM AS-BUILT	Drawings
Sediment thickness (ap		inches				<b>'</b>
Sediment area (approx)	. 10070	1,, X				
				l		
			<i>)</i> .			
No outle	ts U15, 44	1 standing	WATER: NO GOW			
Circle if Applicable:				Signature(s):		
MS/MSD Duplicate ID No.:			F74			



# Tetra Tech, Inc. SOIL & SEDIMENT SAMPLE LOG SHEET

Project Site Nar Project No.: [] Surface So [] Subsurface [X] Sediment [] Other: [] QA Sample	oil e Soil t	Lockheed Martin Midd 112IC05220 Storm S		Sampled E C.O.C. No Type of Sa [] Low C	SD-/74 By: F. Kolberg	p-2
GRAB SAMPLE DAT	THE RESERVE THE PERSON NAMED IN					
Date:	12/2013	Depth Interval	Color		(Sand, Silt, Clay, Moi	
Time:	1130	0"-2"	01.4	VE-F SAN	10 +GILT, 10 are	e, wet
Method:  Monitor Reading (ppn	RAB	0-0	Black	clayey		
COMPOSITE SAMPI	THE RESIDENCE OF THE PARTY OF T	otalia ka kataman maka				
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Moi	eture etc.)
Dale.	Time	Deptil interval	Coloi	Description	(Sand, Sitt, Clay, Wor	sture, etc.,
Method:			LICORTILI persone	3000	500	100
		70.70			20100000 accumum	
Monitor Readings						
(Range in ppm):		No.			2-4000000000000000000000000000000000000	M. Olivatina Works
N/A			senswe			
		-20				
SAMPLE COLLECTI		NON:				
	Analysis		Container Req		Collected	Other
PAHs, PCBs, Moistur	re content		1 - 8-oz wide mouth g	le mouth glass Yes No		
12.12.				-		
eve-						
OBSERVATIONS / N	OTES:	ЭКУКИН ОКИМЕНТИКИ		IMAP:	Diversity in the Committee of	rijessi mir ilijelim
Circle one MRC or M	NAME OF TAXABLE PARTY.					40 1 To 100
Outfall Number:	00	a			Jaki	LEON +TOWERS
Manhole or CB Numb				Lot	6 at MRC	
Manhole or Catch Ba				<b>1</b>	<del></del>	
Material: Conc.	Brick/mortar			1 (	DarkHead Pol	<b>v</b> .
Dimensions: 28	" WIAM.	9.3 Deep		نبر   [		/
Sediment thickness (a		inches		1 151	1 444	4-1
Sediment area (appro	יאסל עוני	ole ;		chesa paghe	10 10 184	H I
		No 11316	e inicks/putlets	1 12		1
		Strange.				
			1	151	Lot 6	,
1		1	1	Paric	Lot 6  Open 41A1  Lot	n
		1	)		last 1	2
				1-0		
<u> </u>		2 61 1	ha	Perul		
Circle if Applicable:	normanianiani	313 51	handing water	Signature(s):	- 24 35 800 8 400 8	
MS/MSD	Duplicate ID	No.		July India (a):		
INIO/INIOU	Duplicate ID	· 11U		11/1		



Project Site Name: Project No.:    Lockheed Martin Middle			Sample Sample Sample C.O.C. N	Location: SD- 155 By: F. Kolberg	b-1	
[X] Sedimen [] Other: [] QA Sampl	t				Sample: Concentration Concentration	
GRAB SAMPLE DA	ΓA:					
Date:	12/ <b>20</b> /2013	Depth Interval	Color		on (Sand, Silt, Clay, Mois	
	12:00 RAB	۵"-۱"	Blacky	M-C S	and ofine grow	u)
Monitor Reading (ppi						and core to the analysis.
COMPOSITE SAMP	LE DATA:					
Date:	Time	Depth Interval	Color	Description	on (Sand, Silt, Clay, Mois	sture, etc.)
Method:					- 10 m	
Monitor Readings		5 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 -			11-11-11	
(Range in ppm):		SC 1000000000000000000000000000000000000			<del></del>	
N/A		8		<u> </u>	200 5 - 21)	
IN/A			· · · · · · · · · · · · · · · · · · ·			
SAMPLE COLLECT	ION INFORMA	TION:				
	Analysis	Silver .	Container Requ		Collected	Other
PAHs, PCBs, Moistu	re content		1 - 8-oz wide mouth glass Yes No			
	7-					ļ
	S		-			<del> </del>
		7000 1			17	
OBSERVATIONS / N	IOTES:		enestamenen mandarietako	MAP:	alitykazuti ir szogornato egyis	
Circle one MRC or N				<b>†</b>		
Outfall Number:	<b>D9</b>			MRC	<b>—</b> —	
Manhole or CB Numi	oer: CB-	88 (uce si	buly)	BILG	1	4
Manhole or Catch Ba			//	ر م	}	ነት
Material: Conc.					ľ	
		1.4 Deep		1	1	
Sediment thickness (	approx): / 1	inches			P.	
Sediment area (appr	ox): 4' x 2' ;	KI" (10050)			t	
$ 2"$ $Rep \rightarrow \rightarrow$				Ramp	30 0147	
Circle if Applicable:				Signature(s)	: 4	
MS/MSD Duplicate ID No.:					FAS	



Project Site Nam Project No.:  [] Surface Soi [] Subsurface [X] Sediment [] Other: [] QA Sample	I Soil Type:	Lockheed Martin Midd 112IC05220 Storm S			ocation: SD- By: F. Kolberg	0-1
Date:	12/ <b>}</b> <i>D</i> /2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture etc.)
	120	Doptil litter var	33.51	- a	and the same of	1 ,
	GRAD	0"-1"	Black	1 -M 34N	o w/ some fin	(5,10054,
Monitor Reading (ppm		,	שלייים		wet	
COMPOSITE SAMPL						
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
Method:						
Monitor Readings						
(Range in ppm):						
N/A						
SAMPLE COLLECTIO	N INFORMA	TION:	natausisesi iski kili jinda		on Warrier by the Mark	Seausilladistusiin
	Analysis		Container Requ	uirements	Collected	Other
PAHs, PCBs, Moisture	content		1 - 8-oz wide mouth g	lass	(Yes) No	
51 11						
OBOSEDVATIONO VAIO	TEO.	All the control of th		IMAD.	main anns a de la company de la company de la company de la company de la company de la company de la company	
OBSERVATIONS / NO	and a supposition of the supposi			MAP:	•	
Circle one MRO or MS Outfall Number:		3	1043	•	}	
Outfall Number: O	U 7 - C B	1076 / 1000	or show	1		
Manhole or Catch Bas		DO IN CCORP	er study)			1
Material: Conc.		•••		]	5D -186-D. T-C807A	′′
Dimensions: 20		x2.9' deep		1 1	7 2-28011	
Sediment thickness (a	pprox):	inches		-	T- nach	
Sediment area (approx	): 20th 2	8"X1"		CB-81	9 MHOON	
				200-	-0	
	$\neg$ $\rightarrow$	7		<b></b>		
	7			Rang		l
1 / /			}	10009		
1 P	7	(	)	1 (		
	13 0	Lep		1		
				}		
Circle if Applicable:		HARARARAKAN SAMARAKAN		Signature(s):		
MS/MSD	Duplicate II	) No.:		1111		
				ry/		



## **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nam Project No.: [] Surface Soi [] Subsurface [X] Sediment [] Other: [] QA Sample	l Soil	Lockheed Martin Midd 112IC05220 Storm S		Sampled C.O.C. No	By: F. Kolbe	
GRAB SAMPLE DATA	A:					
Date:	12/20/2013	Depth Interval	Color	Description	(Sand, Silt, Clay, I	Moisture, etc.)
Time: 1245				E-M CAN	d w/some :	fact Oach
Method: GAA		0"-1"	DEFTAY	100	1 de 1 30 m.c	ווויטק לפיייי
Monitor Reading (ppm	):		4 PINWN	(005c,1	NOI ST	y marca
COMPOSITE SAMPL	E DATA:					
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, I	Moisture, etc.)
Method:						1
Monitor Readings						*****
					, e	
(Range in ppm):	<b>—</b>	1				
N/A	_				www.manage.edu.com.com	
1		ļ				
				A		
SAMPLE COLLECTION	ON INFORMA	ITION:				
	Analysis	M	Container Requ		Collected	Other
PAHs, PCBs, Moisture	content		1 - 8-oz wide mouth glass (Yes) No			
OBSERVATIONS / NO	TEC:			MAP:		
THE RESERVE OF THE PERSON NAMED IN COLUMN 1						
Circle one: MRC or MS Outfall Number: Ot				J Aspha	- 10t)	
Manhala or CB Number	CA.L	1 ( - + 1)		- ASHADI	il .	., I
Manhole or CB Number	in Construction	( DM.S.WC)	•	¥ /25,	e a se soute (	catch
	Brick/mortar				concrete culvert	BASIA N
Dimensions:	Diicivinoriai	alla.			V	(B-4 (IM)
Sediment thickness (a	pprox): 1"	inches				1
Sediment area (approx					7000	
3.17.2.					::: <b>@</b> :/\	
					SD-	187
		1	)		di mental	
		(	)	""		la calca
				No sedim	est in catch.	6431A)
				collected	fromsedinu	nt Accumulated
				in which	+ upstrann of	basia; at Accumulated 5 CB-4
Circle if Applicable:				Signature(s):		
MS/MSD	Duplicate I	D No.:			11/	
		<del></del>		19	X	



## **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nam Project No.: [] Surface Soi [] Subsurface [X] Sediment [] Other: [] QA Sample	il Soil Type:	Lockheed Martin Mido 112IC05220 Storm S			bocation: SD-/88 By: F. Kolberg	) <b>-</b>
GRAB SAMPLE DATA	V 1000 000 000 000 000 000 000 000 000 0	Donath Internet	Color	Decaring ===	(Cand Cile Olari Ba-1	atura ata\
Date: Time: /25°	12/2013	Depth Interval	Color	-1/2	(Sand, Silt, Clay, Mois	
Method: LPAD		0'-1"	DIL brown	FRM SAM	io, trace fines, p	oorly
Monitor Reading (ppm		0-1	and black	grades	louse, moist	
COMPOSITE SAMPL	CORP. THE PERSON NAMED IN COLUMN 2 IN CO.	turiohidikemedinike	nabeanoumagiezotuat	acemanieninesinaesinae	ien viedokturentekontal etetroareita	attacimo a tratue
Date:	Time	Depth interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
Method:						27
Monitor Readings (Range in ppm): N/A						
SAMPLE COLLECTION INFORMATION:  Analysis  PAHs, PCBs, Moisture content			Container Req 1 - 8-oz wide mouth ç		Collected (Yes) No	Other
Dimensions: Sediment thickness (a	SA  er: CB - 8  in Construction  Brick/mortar  pprox):	Couts.le -up	staga)	MAP:  CAtch bas,  Sample v  E Asihari Lot 1	n-we sediment of compositions	Collected Catch Basia D-8 (IRM)
Sediment area (approx):  Circle if Applicable:				concrete culvert	Accomplated sodiment	
MS/MSD	Duplicate ID	No.:			1/	
				1 1/	7	



Project Site Nam Project No.: [] Surface So [] Subsurface [X] Sediment [] Other: [] QA Sample	il Soil	Lockheed Martin Middle River Complex 112IC05220 Storm Sewer Sediment		Sample ID No.: Sample Location: Sampled By: C.O.C. No.:  Type of Sample: [] Low Concentration [] High Concentration		
GRAB SAMPLE DAT	A:					
Date:	12/30/2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
Time: 125			DKGGATANI	F-M SA	ud; little times, f	poorly
Method: GAA		0"-1"	Dk gray and Brown	graded	louse, moist	,
Monitor Reading (ppm	A RESIDENCE		13/own	100000	100000,70000	
COMPOSITE SAMPL	E DATA:					
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
1.500			4881 556		5 T F 1 S F 5 S F	
Method:						
Monitor Readings					**********	
_				-		
(Range in ppm):					770	3
N/A						
Ĭ				2		
SAMPLE COLLECTION	ON INFORMA	TION:				
	Analysis		Container Requ	uirements	Collected	Other
PAHs, PCBs, Moisture	e content	LO HORO.	1 - 8-oz wide mouth g	glass (Yes) No		
	59000000000000			E - 30		
		Production of the second				
ODCEDVATIONS / N	OTEC.		production and the control of the co	IWAD.		A CONTRACTOR OF THE PARTY OF TH
OBSERVATIONS / NO				MAP:		
Circle one: MRC or M						
Outfall Number: 6 ( Manhole or CB Number)	CAL	1. L. W.		Asphant Lot concrete	المريب	ا مد
Mannole or CB Number	er: UD-V	(ONTSIAL)	· . V	رد جوسا	culver CA	The word
Manhole or Catch Bas				( one Mile	10-199 B	Pro (Iran)
	Brick/mortar	N/A		, ,	40-11.	5
Dimensions: Sediment thickness (a	innrov). I	inches		1 7		
Sediment area (appro		IIICIICS				
(4)				· ~~	$\sim$	
				-> ( ·		
				\mathrew \	\~~	
0		/	\	<del></del>		
]		(	} ~	1 A.J	to 1 contract	
				Macan	atch sediment.	
				No sedimu	it in catch bas	11 st. Sample
		_		collected En	M ACCUMULATED LE	-liment
Circle if Applicable:				Signature(s):		
MS/MSD	Duplicate II	D No :		111	CB-6 /Upsen	יארייאין
1110/11100	Dupiloats II				C0-0	I
				' 7 '/		



Project Site Nam Project No.:  [] Surface Soi [] Subsurface [X] Sediment [] Other: [] QA Sample	l Soil	Lockheed Martin Midd 112IC05220 Storm S			Docation: SD-190 By: F. Kolberg	0-1
GRAB SAMPLE DATA	Δ.					
Date:	12/2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
Time:	1445	Dopair into rui				111011111111111111111111111111111111111
	PAB	יו <sub>ו-</sub> ייט ו'	Duck dury +	1-M 3AA	vd, with some worly graded, w	TIA48
Monitor Reading (ppm		1	PLONU	loose pu	orly graded, w	d
COMPOSITE SAMPL	THE RESIDENCE					
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture, etc.)
Method:						
Monitor Readings		-	****			***
_			William A			
(Range in ppm):						
N/A						
					***	
SAMPLE COLLECTION	ON INFORMA	ATION:				ALTHUM DELECTION
	Analysis	1.00	Container Requ	uirements	Collected	Other
PAHs, PCBs, Moisture	content	(i	1 - 8-oz wide mouth glass Yes No		Yes No	
		44.74			and the state of t	
72-1-		200				
				na a Barrier de la contraction del la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contraction de la contractio		010139813801801111111
OBSERVATIONS / NO				MAP:		
Circle one MRO or MS				catch b	AFIN EAST side	. બુ
Outfall Number: Ot		121,000,000		نم لدائده	الساسي	244
Manhole or CB Number				BOULEUM	g. C worth of	
Manhole or Catch Basi Material: Conc.		on:		y drea f	or wave techn	elos ica
		3.3' Deep		\		A I
Dimensions: 23"	21	inches		<b>\</b>	jet fis	kter 4
Sediment thickness (approx		and the second second			Carly Je	N/
Codimont area (approx	<u>, , , , , , , , , , , , , , , , , , , </u>			1	(III) ROA	ا ۱۲ کم
	<del></del>		_	MPC /	CA RAM	, ' I
				Blg,	A KRI	cessed
1	ł	/	\	1" <i>)</i> \		ecusys 2
			)	(	50-140	NE Area 8 Bld C
					1	s aid c
				^		1
			_		1	
Circle if Applicable:				Signature(s):		
MS/MSD	Duplicate I	D No :	water-1997	٠	14	
1110/11100	Jupileate I	2		FJ.	<b>§</b>	



Project Site Nar Project No.: [] Surface So [] Subsurface [X] Sediment [] Other: [] QA Sample	oil e Soil t	Lockheed Martin Midd			By: SD-191 F. Kolberg	)- <u> </u>
GRAB SAMPLE DAT						
Date:	12/20/2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
Time:  Method:  Monitor Reading (ppr	1450 LAB m): IVIA	0"-1"	Black/ Brown		who wisome fin	ies, 1005 c
COMPOSITE SAMP	Name and Address of the Owner, where the		District Kinnes (Co. 1984)			
Date:	Time	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	sture, etc.)
Method:					20000	7, 12
Monitor Readings	<del>                                     </del>					
(Range in ppm):						
N/A					manufi we	
11/7	<b>-</b>					P
PAHs, PCBs, Moistur	Analysis re content		Container Req		Yes No	Other
OBSERVATIONS / N				MAP:		
Circle one MRC or M				T - 4 1 1		rat 1 mars
Outfall Number:	707				as in located be	
Manhole or CB Numb				TWC B	11de c entance	ec near
Manhole or Catch Ba		on:		neck 8	and Alcui	not A
Material: Conc.  Dimensions: 24		x 2.9' Deep		1	and A/C united	4
Sediment thickness (a		inches		/ CB N	clas ALC DW	W
Sediment area (appro	ox): 100 %			1 / 20	cassed the step	15
	No	pipe visible -	c8 flooded	1 \ 1	TO A LAND	Road
		.,		TYPE 2:M/	7 1	Jet Fiskley
				Blag c Y	30-MI	,
		(		1 2.27 2	W) /-	RAMY
		\	/	1111	V-A/c	
				1 / / /	(E) outed)	/
		Standing water		1 / /	/ / \	/
Circle if Applicable:				Signature(s):		
MS/MSD	Duplicate II	D No.:		FOR	<i>*</i> ,	
INVESTIGATION OF THE PROPERTY				177		



# Tetra Tech, Inc. SOIL & SEDIMENT SAMPLE LOG SHEET

Project Site Nam Project No.: [] Surface Soi [] Subsurface [X] Sediment [] Other: [] QA Sample	l Soil Type:	Lockheed Martin Midd	<del></del>		y: SD-192 F. Kolberg	-0-1
GRAB SAMPLE DATA						
Date:	12/20/2013	Depth Interval	Color	Description (	Sand, Silt, Clay, Moi	sture, etc.)
	00	0"-1"	مستواطام	ـــ	1 .11 4	
Method: Gr		0 -1	Light brown	F-c sand, w/trace fines Pourly graded, louse, wet		
Monitor Reading (ppm		CMAR market by the control of the co		Poorly 9	raded 1 louse	IWET
COMPOSITE SAMPL	E DATA:	· · · · · · · · · · · · · · · · · · ·	·			
Date:	Time	Depth Interval	Color	Description (	Sand, Silt, Clay, Moi	sture, etc.)
	centros en esto-en en el			r.		
Method:						
		1 2.11			₹	
					1	
Monitor Readings						
(Range in ppm):						
N/A			8			
SAMPLE COLLECTION	N INFORMA	TION:				
	Analysis		Container Requ	irements	Collected	Other
PAHs, PCBs, Moisture			1 - 8-oz wide mouth glass Yes No			Other
r Ai is, r Obs, Moisture	Content		1 o de wide mount glass			1
				<del> </del>		
nominos are				-		
OBSERVATIONS / NO	TEC.	Rundedtonandinuschilinis		MAP:		
		recuss ratssame autometric		MAP.	ENDRE CHRUMULARI MULURUS	DATE OF THE PROPERTY OF
Circle one: MRC or MS		1 .4 14 - 1	Conld	1 M	1	A I
Outfall Number:	MASTICAM (	AND WHOL	OUTA		1	
Manhole or CB Number				Alder	I	1/
Manhole or Catch Bas		n:		Alder	1	ן אי
Material: Conc	the state of the s	21 1			1	•
Dimensions: 7' x		3' deep			- 1	
Sediment thickness (a	pprox):	inches	1 -		> 1	
Sediment area (approx	9: 5070	- MOSTLY COM	bus	Baire	E (0)92	
			_		_ AU	1601
				11/7	WROF	1
à .	Į			1	WRDF	1001 A
		1	1	8	- 1	
		(	1	3	• 1	1
				-2	· 1	
			/	13	1	
•				]	ı	
Circle if Applicable:				Signature(s):		
M2/M2D 1	Duplicate ID	) No.:		City	1 Y	I
MS/MSD	Duplicate ID	No.:		FI	X	



## **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Na Project No.: [] Surface S [] Subsurfact [X] Sedimet [] Other: [] QA Samp	Soil ce Soil nt	Lockheed Martin Midd 112IC05220 Storm 9		Sample ID No.: Sample Location: Sampled By: C.O.C. No.:  Type of Sample: Y Low Concentration [] High Concentration
GRAB SAMPLE DA	ATA:			
Date:	12 <b>20</b> /2013	Depth Interval	Color	Description (Sand, Silt, Clay, Moisture, etc.)
Time:	1610	• •	Brown	
	GRAB	0"-1"	Diemy	F-c sand w/ some fines, lover, wet, graded
Monitor Reading (p	THE RESERVE OF THE PERSON NAMED IN		Malaria Nicolana	Wet, granen
COMPOSITE SAM	PLE DATA:			
Date:	Time	Depth Interval	Color	Description (Sand, Silt, Clay, Moisture, etc.)
Method:		100		
Monitor Readings				
(Range in ppm):	-	19.25		(A) (A) (A) (A) (A) (A) (A) (A) (A) (A)
		3/2		
N/A				
PAHs, PCBs, Moist	are content	083	1 - 8-oz wide mouth g	glass Yes No
		Para-		
OBSERVATIONS /	NOTES:			MAP:
Circle one: MRC of	and the second			
		ed ditches f	ram MSA	sample collect in open chandal
Manhole or CB Nun	nber: NA	of ditches f		The state of the s
Manhole or Catch E				(Small ditch) pear headwall
Material: Conc	Brick/mortar			يومان ۽ مسم پريانون
Dimensions:				Across ton 9000 wilson
Sediment thickness		inches		(Small ditch) year headwall Across from 900B wilson point Rd
Sediment area (app				Headush
Circle if Applicable		D.No.:		Signature(s):
MS/MSD	Duplicate II	אס.:		F! A



# **SOIL & SEDIMENT SAMPLE LOG SHEET**

Project Site Nam Project No.: [] Surface Soi [] Subsurface [X] Sediment [] Other: [] QA Sample	- I Soil	Lockheed Martin Midd 112IC05220 Storm S			SD-194  By: F. Kolberg	D-I
GRAB SAMPLE DATA	A:					
Date:	12/ <b>}0</b> /2013	Depth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture, etc.)
Time:	AB	0'-1"	LT Brown		d w/some f.n	
COMPOSITE SAMPL			inimaran kanazarat Patri			
	The second second	Donth Interval	Color	Description	(Sand, Silt, Clay, Mois	ture. etc.)
Date:	Time	Depth Interval		Description	(Ourie, Only Olay) more	
Method:						
Monitor Readings		-				12 23 1
(Range in ppm): N/A						
SAMPLE COLLECTION		TION:	L Containe Boo		Collected	Other
	Analysis		Container Rec		Collected No	Other
PAHs, PCBs, Moisture	e content		1 - 8-oz wide mouth	yiass	(Yes) No	
			<del> </del>			
			<u> </u>	725		
OBSERVATIONS / N	OTES:			MAP:		
Circle one: MRC or M	A STREET, SQUARE					
		AM & MSA L	10 OF ODIA	Sample o	collected from . eatch basin in	storm
Manhole or CB Numb	er:	AN PINON -	7401		411	claset
Manhole or Catch Bas		on:		d'am	cateu basia in	371 001
Material: Conc.	Brick/mortar			IN EN,	it of 9008 w fund. enm of 50-199	·ISUN
Dimensions:		31		l as it	0.00	
Sediment thickness (a	approx):	inches		I porta	MUNIC.	
Sediment area (appro	x):			044	- · · · · · · · · · · · · · · · · · · ·	7.1
				50-	-193	
Circle if Applicable:				Signature(s):		
MS/MSD	Duplicate I	D No.:		] FJ	#	

APPENDIX B—DATA VALIDATION REPORTS (ON CD ONLY)	



#### INTERNAL CORRESPONDENCE

TO:

T. APANAVAGE

DATE:

**JANUARY 28, 2014** 

FROM:

MICHELLE L. ALLEN

COPIES:

**DV FILE** 

SUBJECT:

ORGANIC DATA VALIDATION - PAH/PCB

LOCKHEED MARTIN CORPORATION (LMC) - SEDIMENT INVESTIGATION

SDG 180-28384-1

SAMPLES:

32/Sediment/PAH/PCB

SD-163-0-1	SD-164-0-1	SD-165-0-1
SD-166-0-1	SD-167-0-1	SD-168-0-1
SD-169-0-1	SD-170-0-1	SD-171-0-1
SD-172-0-3	SD-173-0-3	SD-174-0-1
SD-175-0-1	SD-176-0-3	SD-177-0-2
SD-178-0-1	SD-179-0-1	SD-180-0-1
SD-181-0-1	SD-182-0-3	SD-183-0-1
SD-184-0-2	SD-185-0-1	SD-186-0-1
SD-187-0-1	SD-188-0-1	SD-189-0-1
SD-190-0-1	SD-191-0-1	SD-192-0-1
SD-193-0-1	SD-194-0-1	

#### **Overview**

The sample set for LMC, SDG 180-28384-1 consisted of thirty-two (32) sediment environmental samples. All thirty-two (32) sediment samples were analyzed for polynuclear aromatic hydrocarbons (PAH) and polychlorinated biphenyls (PCB). No field duplicate sample pair was included in this SDG.

The samples were collected by Tetra Tech, Inc. on December 18-20, 2013 and analyzed by TestAmerica, Inc. All analyses were conducted in accordance with SW-846 Methods 8270D and 8082A analytical and reporting protocols.

The data contained in this SDG were validated with regard to the following parameters: data completeness, holding times/sample preservation, GC/MS tuning, initial/continuing calibrations, laboratory method blank results, surrogate spike recoveries, blank spike results, matrix spike/matrix spike duplicate results, internal standard areas and recoveries, chromatographic resolution, compound identification, and detection limits. Areas of concern are listed below.

#### <u>Major</u>

No major issues were identified.

#### Minor

The Percent Solids (% Solids) was less than 30% in the following samples:

<u>Sample</u> <u>% Solids</u> SD-177-0-2 71.8

The detected and non-detected results reported for the PAH target compounds and PCBs in these samples were qualified as estimated, (J) and (UJ), respectively.

TO: T. APANAVAGE SDG: 180-28384-1

• The internal standard areas for perylene-d12 were above the upper quality control limit in samples SD-169-0-1, SD-170-0-1, SD-171-0-1, SD-172-0-3, SD-173-0-3, SD-174-0-1, SD-175-0-1, SD-176-0-3, SD-177-0-2, SD-178-0-1, and SD-179-0-1. These samples were not re-extracted or reanalyzed. The detected and non-detected results reported for the PAH compounds associated with this internal standard were qualified as estimated, (J) and (UJ), respectively.

 The PAH Matrix Spike/Matrix Spike Duplicate analyses performed on sample SD-185-0-1 had the following noncompliances:

<u>Compound</u>	<u>MS %R</u>	MSD %R	<u>RPD</u>	QC Limits	RPD Limit	Qualifier
Anthracene	131*	59	49*	43-111	35	NA
Benzo(a)anthracene	21*	-89*	42*	45-110	31	J
Benzo(b)fluoranthene	-4*	-98*	34*	37-108	28	J
Benzo(k)fluoranthene	82	26*	31	39-115	42	NA
Benzo(g,h,i)perylene	59	-20*	29*	35-127	21	NA
Benzo(a)pyrene	29*	-74*	40*	42-114	31	J
Chrysene	47	-105*	46*	44-108	31	NA
Fluorene	131*	87	3	43-110	37	NA
Fluoranthene	-58**	-369**	52*	40-118	23	NA
Indeno(1,2,3-cd)pyrene	40	-20*	25	34-130	30	NA
Phenanthrene	181*	-46*	75*	41-107	20	J
Pyrene	24**	-170**	56	39-113	28	NA
Acenaphthene	119*	78	36*	42-104	34	NA

<sup>\* -</sup> outside QC limit

NA - No action

No action was taken in the parent sample if at least one %R was within the quality control limits. No action was taken for noncompliant RPDs because either one or both %Rs were within the quality control limits. In addition, not action was taken for fluoranthene and pyrene because the sample concentrations of these compounds were greater than 4X the concentration spiked in the sample and the MS/MSD samples were analyzed at a 20X dilution. The remaining noncompliances were qualified as indicated in the parent sample.

- The PAH MS/MSD analyses performed on sample SD-163-0-1 (50X dilution) had %Rs for all the target compounds greater than the upper quality control limits with the exception of dibenz(a,h)anthracene and naphthalene. The MS %R for naphthalene was above the upper quality control limit (42-104) but the MSD %R and RPD were acceptable, therefore, no action was taken. In addition, no action was taken for 2-methylnaphthalene because this compound was not detected in the parent sample. The detected results reported above the Reporting Limit (RL) in the parent sample for the remaining PAH compounds were qualified as estimated, (J), due to conflicting directional bias.
- The PCB MS/MSD analyses performed on sample SD-165-0-1 had %Rs for Aroclor 1260 below the lower quality control limit. The detected result reported for this PCB in the parent sample was qualified as estimated, (J).
- Detected results reported below the RL limit but above the Method Detection Limit (MDL) were qualified as estimated, (J).

<sup>\*\* -</sup> outside QC limit, sample concentration > 4X spiked concentration

J - Estimated due to conflicting directional bias

TO: T. APANAVAGE SDG: 180-28384-1

#### <u>Notes</u>

All sample PCB chromatograms were reviewed for evidence of Aroclor mixture dechlorination and/or weathering as similarly encountered in other sediment samples located within the Middle River Complex site. Upon review, some samples contained peaks that could possibly be indicative of PCB dechlorination/weathering but may also be indicative of other non-PCB contaminants. The samples that contain peaks that elute outside of the reported Aroclor mixture but are not similar to previous samples that exhibited evidence of dechlorination are as follows: SD-168-0-1, SD-172-0-3, SD-184-0-2, SD-187-0-1, and SD-189-0-1. No action was taken based on the uncertainty and limitations associated with the sample data.

One sample, SD-182-0-3, does appear to exhibit a pattern similar to previous samples where dechlorination/weathering was suspected. This sample contained the highest concentration of Aroclor 1260 in this data set and displayed three to five peaks (between approximately 6 and 8 minutes) that eluted prior to the primary Aroclor 1260 pattern. It is worth noting that the Aroclor 1260 pattern in this sample was a very good match when compared to the standard Aroclor 1260 pattern. The unidentified peaks in the sample were similar to those encountered in other sediment samples that were suspected of exhibiting dechlorination/weathering but the peak intensities were not considered significant by the data reviewer because the magnitude of the peaks were less than the overall intensity of the Aroclor 1260 pattern. No validation action was taken.

The %Rs for Aroclor 1016 in the MS/MSD analyses of sample SD-185-0-1 were above the upper quality control limit. In addition, the MSD %R was low for Aroclor 1260. No action was taken because Aroclor 1016 was not detected in the parent sample. No action was taken for Aroclor 1260 because the concentration of this PCB in the parent sample was greater than 4X the amount spiked in the MS/MSD samples.

Second column confirmation was not performed on the samples that had PCB detections. However, PCB mixtures can be identified on the basis of chromatographic patterns; therefore, there is no effect on sample identification.

All the PAH and PCB samples were initially analyzed at dilutions. Consequently, the reporting limits of the non-detected PAH and PCBs were elevated and the surrogate spike compounds were not recovered in many samples.

The following samples required further dilution due to PAH compounds exceeding the calibration range of the instrument. The results from the dilutions were used in the data validation.

<u>Sample</u>	<u>Compound</u>	<u>Dilution</u>
SD-190-0-1	Fluoranthene	50X
SD-191-0-1	Fluoranthene	100X
	Phenanthrene	100X
	Pyrene	100X

Non-detected results were reported to the MDL.

TO: T. APANAVAGE SDG: 180-28384-1

#### **Executive Summary**

**Laboratory Performance:** The %Solids exceeded 30% in one sample. Several PAH samples had high internal standard areas for perylene-d12. Second column confirmation was not performed in the PCB fraction.

**Other Factors Affecting Data Quality:** All samples were diluted in the PAH and PCB fractions. The PAH and PCB MS/MSD analyses had noncompliant %Rs. Potential PCB dechlorination was present in one sample. Results below the RL were estimated.

The data for these analyses were reviewed with reference to Region III modifications to U.S. EPA National Functional Guidelines for Organic Data Validation (Sept. 1994) and SW-846 Methods 8270D and 8082A analytical and reporting protocols. 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

Terfa 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

PROJ NO: 05483	NSAMPLE	SD-163-0-1			SD-164-0-1			SD-165-0-1			SD-166-0-1		
SDG: 180-28384-1	LAB_ID	180-28384-1			180-28384-2			180-28384-3			180-28384-4		
FRACTION: PAH	SAMP_DATE	12/18/2013			12/18/2013			12/18/2013			12/18/2013		
MEDIA: SOIL	QC_TYPE	ΣZ			ΣN			NM			NN		:
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG		
	PCT_SOLIDS	76.8			78.8			76.2			80.0		
	DUP_OF							·					
PARAMETER		RESULT	Val	alcd	RESULT	VQL	arcp	RESULT	NOL	arcd	RESULT	VQL	QLCD
2-METHYLNAPHTHALENE		19 U	D		30	7	а	25	25 J	Ь	12	75 U	
ACENAPHTHENE		64	٦	DP	25	7	Ь	47	ſ	Ь	8	80 U	
ACENAPHTHYLENE		69	ſ	DP	150	ſ	Ь	36	ſ	Ь	ō	0 96	
ANTHRACENE		26	ſ	DP	200	ſ	Ь	110	7	Д	88	82 U	
BENZO(A)ANTHRACENE		530	ſ	О	520			830			130	r o	۵
BENZO(A)PYRENE		009	ſ	D	099			920			120	7	Д
BENZO(B)FLUORANTHENE		200	ſ	D	640			1400			13.	130 U	
BENZO(G,H,I)PERYLENE		480 J	ſ	D	760			890			13	130 J	<u>م</u>
BENZO(K)FLUORANTHENE	==	260	J	D	420			290			171	170 U	
CHRYSENE		710 J	J	D	760			1500			180	ر 0	Д.
DIBENZO(A,H)ANTHRACENE	JE JE	110	ſ	DP	130	ſ	Ь	200	7	Д	100	<u></u>	Д
FLUORANTHENE		1400	ſ	О	096			3500			180	ر o	۵
FLUORENE		55	ſ	DP	52	ſ	Д	55	r	Д	110	110 U	-
INDENO(1,2,3-CD)PYRENE		390	ſ	Q	260			790			æ,	86 U	
NAPHTHALENE		19	n		34 J	ſ	Ь	95	ſ	Ь	2	72 U	
PHENANTHRENE		066	ſ	D	370			2800			130	n	
PYRENE		1100	ſ	۵	1100			2500			210	7	Ь

PROJ_NO: 05483	NSAMPLE	SD-167-0-1			SD-168-0-1			SD-169-0-1			SD-170-0-1		
SDG: 180-28384-1	LAB_ID	180-28384-5			180-28384-6			180-28384-7			180-28384-8		
FRACTION: PAH	SAMP_DATE	12/18/2013			12/18/2013			12/18/2013			12/20/2013		
MEDIA: SOIL	QC_TYPE	NN			NN			NM			NN		
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG		
	PCT_SOLIDS	9.77			78.7			97.6			78.5		
	DUP_OF												
PARAMETER		RESULT	VaL	alcd	RESULT	VQL	arcd	RESULT	VQL	arcd	RESULT	VQL	arcd
2-METHYLNAPHTHALENE		39	_		38	Э		17	_			19 U	
ACENAPHTHENE		41	n		110	ſ	Ь	30	30 J	Ь	2	20 U	
ACENAPHTHYLENE		49			48 U	_		22	22 U		5	58 J	Ь
ANTHRACENE		44	ſ	Ь	180	ſ	Ь	19	n		110	٥ ر	Ь
BENZO(A)ANTHRACENE		270	_	۵	470			92	٦	Ь	320	0	
BENZO(A)PYRENE		280	ſ	Ь	390	7	Ь	69	7	NP	250	ا م	Z
BENZO(B)FLUORANTHENE		430			250	٦	Ь	120	ſ	ΝΡ	320	0 ر	z
BENZO(G,H,I)PERYLENE		340 J	ſ	Ь	410	ſ	Ь	81	J	NP	170	٥ ر	NP
BENZO(K)FLUORANTHENE		170	ſ	Д.	530			43	ſ	ΝΡ	2	78 J	NP
CHRYSENE		440			530			120 J	ſ	С	280	0	
DIBENZO(A,H)ANTHRACENE	무	48 U	n		88	J	Ь	21	21 UJ	z	8	32 J	NP
FLUORANTHENE		800			1000			180 J	٦	Ь	930	.0	
FLUORENE		N 29	n		98	ſ	Ь	25	25 U		2	28 U	
INDENO(1,2,3-CD)PYRENE		1 00E	ſ	Ь	370	ſ	Ь	99	r 99	Νb	150	ا ر	NP
NAPHTHALENE		0 7E	n		98	ſ	Ь	16	Π		1	18 U	
PHENANTHRENE		340	ſ	Ь	710			47	47 J	Ь	300	0	
PYRENE		290			830			160	ſ	Ь	480	0	

PROJ_NO: 05483	NSAMPLE	SD-171-0-1			SD-172-0-3			SD-173-0-3			SD-174-0-1		
SDG: 180-28384-1	LAB_ID	180-28384-9			180-28384-10	:		180-28384-11			180-28384-12		
FRACTION: PAH	SAMP_DATE	12/19/2013			12/19/2013			12/19/2013			12/19/2013		
MEDIA: SOIL	QC_TYPE	NN		į	NM			NM			NM		
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG		
	PCT_SOLIDS	33.2			53.1			62.0			83.4		
	DUP_OF												
PARAMETER		RESULT	ΛQL	QLCD	RESULT	VQL	arcp	RESULT	VQL	arcp	RESULT	VQL	QLCD
2-METHYLNAPHTHALENE		14	_	۵	90	ſ	Ь	55	ſ	Ь	47	J	Ь
ACENAPHTHENE		28	_	Ь	920			150	ſ	Ь	230		
ACENAPHTHYLENE		71			160	ſ	Д	430			140	ſ	Д.
ANTHRACENE		160			1100			580			610		
BENZO(A)ANTHRACENE		620			4600			1700			2200		
BENZO(A)PYRENE		089	_	z	4200 J	ſ	Z	1900	J	Z	1900	ſ	z
BENZO(B)FLUORANTHENE	=	720	ſ	z	5300 J	ſ	Z	2100	<b>ا</b>	z	2500	7	z
BENZO(G,H,I)PERYLENE		C 089	_	z	3300	J	Z	1800	J	z	1300	7	z
BENZO(K)FLUORANTHENË	711	250	_	z	1800	ſ	Z	630	r	z	780	7	z
CHRYSENE		720			2000			2000			2400		
DIBENZO(A, H)ANTHRACENE	<b>LE</b>	150	7	z	870	J	Z	390	ſ	z	360	7	z
FLUORANTHENE		910			0096			3000			5800		
FLUORENE		35	_	۵	460			160	ر ا	Ь	250		
INDENO(1,2,3-CD)PYRENE		930	7	z	3100	ſ	Z	1500	ſ	z	1300	7	z
NAPHTHALENE		24	_	<u>а</u>	88	ſ	Ь	94	ſ	Ь	160	ſ	Д.
PHENANTHRENE		430			5000			1100			3800		
PYRENE		1100			7400			3400			4000		

PROJ_NO: 05483	NSAMPLE	SD-175-0-1			SD-176-0-3			SD-177-0-2		SD-178-0-1	
SDG: 180-28384-1	LAB_ID	180-28384-13			180-28384-14	_		180-28384-15		180-28384-16	
FRACTION: PAH	SAMP_DATE	12/19/2013			12/19/2013			12/19/2013		12/19/2013	
MEDIA: SOIL	QC_TYPE	ΣZ			ΝN			NN		NM	
	UNITS	UG/KG			UG/KG			UG/KG		UG/KG	
	PCT_SOLIDS	70.0			56.7			28.2		78.6	
	DUP_OF										
PARAMETER		RESULT	VaL	arcd	RESULT	VQL	arcp	RESULT VOL	arcd	RESULT VOL OL	alcd
2-METHYLNAPHTHALENE		210	_	Ь	30	7	Ь	300	ΡΥ	19 U	
ACENAPHTHENE		1500			130	7	۵	2200 J	٨	20 U	
ACENAPHTHYLENE		310			200	7	ď	510 J	ΡY	48 J P	
ANTHRACENE		2100			390	[(		3300 J		30 J P	
BENZO(A)ANTHRACENE		9220			1600			18000	<b>&gt;</b>	140 J P	
BENZO(A)PYRENE		5400	ſ	z	1800 J	٦ ر	z	21000 J	N	180 J NP	
BENZO(B)FLUORANTHENE	ш	0009	ſ	z	2700	r l	z	27000 J	Ν	200 J NP	
BENZO(G,H,I)PERYLENE		3700	7	z	1500	ر ا	Z	15000 J	ΛN	140 J NP	
BENZO(K)FLUORANTHENE	Е	2500	ſ	z	710	٦ ر	Z	10000	Ϋ́	86 J NP	
CHRYSENE		6100			2200	_		26000 J	<b>\</b>	190 J P	
DIBENZO(A, H)ANTHRACENE	NE	1100	ſ	z	320	ſ	z	3500 J	Ν	32 J NP	
FLUORANTHENE		13000			4200			C 00089	٨	260	
FLUORENE		1000			160	r (	Ь	2200 J	٨	28 U	
INDENO(1,2,3-CD)PYRENE	=	r   009£	ſ	Z	1400 J	ر (	z	14000 J	Σ	110 J NP	
NAPHTHALENE		480			51	ר	Ь	120 J	ΡY	18 U	
PHENANTHRENE		9000			1700			39000 J	<b>\</b>	81 J	
PYRENE		10000			3000			ל 45000	Υ	210	

PRO.1 NO: 05483	NSAMPI F	SD-179-0-1			SD-180-0-1			SD-181-0-1		SD-182-0-3	
SDG: 180-28384-1	LAB_ID	180-28384-17			180-28384-18			180-28384-19		180-28384-20	
FRACTION: PAH	SAMP_DATE	12/19/2013			12/19/2013			12/20/2013		12/20/2013	
MEDIA: SOIL	QC_TYPE	NA NA			ΣN			MN		NM	
	UNITS	UG/KG			UG/KG			UG/KG		UG/KG	
	PCT_SOLIDS	2.69			78.7			57.5		47.9	
	DUP_OF										
PARAMETER		RESULT	VQL	arcp	RESULT	VQL	arcp	RESULT VQL	arcd	RESULT VOL	QLCD
2-METHYLNAPHTHALENE		2.5	<u></u>	۵	38	n		440 J	Ь	240 J	Ь
ACENAPHTHENE		18			70	7	Ь	3900		1600	
ACENAPHTHYLENE		14			92	ſ	۵	350 J	Ь	230 J	Ь
ANTHRACENE		41			390	7	۵	7100		3000	
BENZO(A)ANTHRACENE		120			2600	_		20000		10000	
BENZO(A)PYRENE		200	<u></u>	z	2900			17000		9200	
BENZO(B)FLUORANTHENE	ш	290 J	ſ	Z	4300			21000		13000	
BENZO(G,H,I)PERYLENE		150 J	ſ	z	2600			14000		9200	
BENZO(K)FLUORANTHENE	Е	120 J	۲	z	1900			8800		3500	
CHRYSENE		230			3700			20000		11000	
DIBENZO(A,H)ANTHRACENE	JN.	32	ſ;	Z	490			3100		2200	
FLUORANTHENE		410			6200	_		35000		18000	:
FLUORENE		13	1		110	ſ	Ь	4100		1200	
INDENO(1,2,3-CD)PYRENE		150	ſ	z	2200			12000		7300	
NAPHTHALENE		4.7	ſ,	Ь	36	. n		640		450 J	Ь
PHENANTHRENE		150			2600			28000		12000	
PYRENE		290			5900			37000		18000	

PROJ_NO: 05483	NSAMPLE	SD-183-0-1			SD-184-0-2			SD-185-0-1			SD-186-0-1			
SDG: 180-28384-1	LAB_ID	180-28384-21			180-28384-22			180-28384-23			180-28384-24			
FRACTION: PAH	SAMP_DATE	12/20/2013			12/20/2013			12/20/2013			12/20/2013			
MEDIA: SOIL	QC_TYPE	ΣZ			ΣN			NN			NM			
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG			
	PCT_SOLIDS	30.1			67.9			74.6			74.2			
	DUP_OF													
PARAMETER		RESULT	VaL	arcd	RESULT	VaL	arcp	RESULT	NOF (	OLCD	RESULT	VQL	arcd	
2-METHYLNAPHTHALENE		280			170			21 J		Ь	32	32 U		-
ACENAPHTHENE		3000			920			53 J		Ь	35	35 U		
ACENAPHTHYLENE		320			260			94 J		Ь	50 J	7	Ъ	
ANTHRACENE		2600			2200			230			96	7	а	
BENZO(A)ANTHRACENE		15000			4500			1300		٥	280	7	Ь	
BENZO(A)PYRENE		13000			3800			1200 J		٥	290 J	7	Ь	
BENZO(B)FLUORANTHENE	Ш	15000			4700			1500	)	D	540	i		
BENZO(G,H,I)PERYLENE		11000			3400			1100			350	7	Ъ	
BENZO(K)FLUORANTHENE	Ш	6200			1700			570			150	7	Ъ	
CHRYSENE		16000	_		4900			1600			520			
DIBENZO(A,H)ANTHRACENE	JN.	3100			930			270			29	7	Ь	
FLUORANTHENE		33000	_		11000			3600			1100			
FLUORENE		2400			1100			87	7	Д	47	D	-	
INDENO(1,2,3-CD)PYRENE	Ш	9200			3000			1000			300	7	Д	
NAPHTHALENE		480			340			43 J		_	31	D		
PHENANTHRENE		17000	_		6400			1100		0	400			
PYRENE		19000			6100			1900			590			

PROJ_NO: 05483	NSAMPLE	SD-187-0-1			SD-188-0-1			SD-189-0-1			SD-190-0-1		
SDG: 180-28384-1	LAB_ID	180-28384-25			180-28384-26			180-28384-27			180-28384-28		
FRACTION: PAH	SAMP_DATE	12/20/2013			12/20/2013			12/20/2013			12/20/2013		
MEDIA: SOIL	QC_TYPE	ΣN			NN			NM			Z		
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG		
	PCT_SOLIDS	79.7	i		88.3			90.6			8.69		
-	DUP_OF												
PARAMETER		RESULT	VaL	arcp	RESULT	VQL	arcp	RESULT	VQL	alcd	RESULT	VaL	QLCD
2-METHYLNAPHTHALENE		48	48 J	۵	27	n		26 U	n		530		
ACENAPHTHENE		48	48 U		29	n		130	ſ	Ь	5600		
ACENAPHTHYLENE		190	<b>¬</b>	Ь	120 J	7	Ь	160 J	7	Ь	870		
ANTHRACENE		260	_	Д	140 J	ſ	Ь	590			18000		
BENZO(A)ANTHRACENE		1100	_		470			3000			49000		
BENZO(A)PYRENE		1400			620			3300			43000		
BENZO(B)FLUORANTHENE	ш	1800	_		800			3900			47000		
BENZO(G,H,I)PERYLENE		1600			069			2800			33000		
BENZO(K)FLUORANTHENE	Ш	720			250 J	ſ	Ь	1500			19000		
CHRYSENE		1700			710			3700			55000		
DIBENZO(A,H)ANTHRACENE	NE	340	ſ	Ъ	150 J	7	Ь	089			8200		
FLUORANTHENE		2900	_		1000			8100					
FLUORENE		82	ſ	Ь	43 J	7	Ь	180	7	۵	0099		
INDENO(1,2,3-CD)PYRENE	•••	1300			540			2600			30000		
NAPHTHALENE		43	43 U		26	n n		P 99	7	Д	120	7	Ь
PHENANTHRENE		850	-		260	7	Ь	2700			61000		
PYRENE		1900	_		710			4500			74000		

PROJ_NO: 05483	NSAMPLE	SD-190-0-1DL			SD-191-0-1			SD-191-0-1DL			SD-192-0-1		
SDG: 180-28384-1	LAB_ID	180-28384-28			180-28384-29			180-28384-29			180-28384-30		
FRACTION: PAH	SAMP_DATE	12/20/2013			12/20/2013			12/20/2013			12/20/2013		
MEDIA: SOIL	QC_TYPE	NN			NM			NM			NM		
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG		
	PCT_SOLIDS	8.69			6.69			6.69			76.2		
	DUP_OF												
PARAMETER		RESULT	VQL	QLCD	RESULT	VQL	QLCD	RESULT	VQL	arcd	RESULT	VQL	QLCD
2-METHYLNAPHTHALENE					230						7.8	n s	
ACENAPHTHENE					2000			:			20	l J	Ь
ACENAPHTHYLENE					440						89		
ANTHRACENE					4200						88		
BENZO(A)ANTHRACENE					13000						130		
BENZO(A)PYRENE					13000						150		
BENZO(B)FLUORANTHENE	111				16000						170		
BENZO(G,H,I)PERYLENE					9200						120		
BENZO(K)FLUORANTHENE	III				5600						64	7	<u>م</u>
CHRYSENE					17000						190		
DIBENZO(A,H)ANTHRACENE	NE				2300						29	7	Ь
FLUORANTHENE		150000						67000			360		
FLUORENE					2400						34	7	<u>_</u>
INDENO(1,2,3-CD)PYRENE					8500						66		
NAPHTHALENE					87	J	Ь				7.5	i U	
PHENANTHRENE				-				42000			170		
PYRENE								49000		:	240		

PROJ_NO: 05483	NSAMPLE	SD-193-0-1			SD-194-0-1		
SDG: 180-28384-1	LAB_ID	180-28384-31			180-28384-32		
FRACTION: PAH	SAMP_DATE	12/20/2013			12/20/2013		
MEDIA: SOIL	QC_TYPE	ΣN			ΝN		
	UNITS	UG/KG			UG/KG		
	PCT_SOLIDS	70.8			84.3		
	DUP_OF						
PARAMETER		RESULT	VQL	arce	RESULT	NOL	arcd
2-METHYLNAPHTHALENE		8.4 U	n		7.1	n	
ACENAPHTHENE		13	ſ	Ь	6	J	Ь
ACENAPHTHYLENE		31	ſ	Ь	48	ſ	Ь
ANTHRACENE		54	ſ	Ь	43 J	ſ	Ь
BENZO(A)ANTHRACENE		160			100		•
BENZO(A)PYRENE		170			130		
BENZO(B)FLUORANTHENE	==	200			140		
BENZO(G,H,I)PERYLENE		130			94		
BENZO(K)FLUORANTHENE		28	٦	Ь	44	ſ	Ь
CHRYSENE		200			150		
DIBENZO(A,H)ANTHRACENE	JE JE JE JE JE JE JE JE JE JE JE JE JE J	33	ſ	Ь	26	ſ	Ь
FLUORANTHENE		440			240		
FLUORENE		23 J		Д	15	ſ	Ь
INDENO(1,2,3-CD)PYRENE		110			62	ſ	Ь
NAPHTHALENE		8.1	n		8.9	n	
PHENANTHRENE		160			81		
PYRENE		230			160		

PROJ_NO: 05483	NSAMPLE	SD-163-0-1			SD-164-0-1			SD-165-0-1	•		SD-166-0-1			
SDG: 180-28384-1	LAB_ID	180-28384-1			180-28384-2			180-28384-3	~		180-28384-4			
FRACTION: PCB	SAMP_DATE	12/18/2013			12/18/2013			12/18/2013			12/18/2013			
MEDIA: SOIL	QC_TYPE	ΣZ			ΣN			ΣN			ΣN			
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG			
	PCT_SOLIDS 76.8	76.8			78.8			76.2			80.0			
-	DUP_OF													
PARAMETER		RESULT	VaL	arcp	RESULT	VaL	arcd	RESULT	ΛQΓ	OLCD	RESULT	VQL	arcd	
AROCLOR-1016		0.4	4 U		0.39	D		0.41	11 U		0.39	Ωle		
AROCLOR-1221		0.52	2 U		0.5	D		0.5	0.52 U		0.5	)    -		
AROCLOR-1232		0.46	9 <u>۱</u>		0.45	D		7:0	0.47 U		0.45	0 9		
AROCLOR-1242		0.44	4 U		0.43	) ]		7.0	0.44 U		0.42	O 3		
AROCLOR-1248		0.26	9 N		0.25	D		7.0	0.26 U		0.25	. n		
AROCLOR-1254		0.38	8 U		0.38			0.5	0.39 U		0.37	u/		
AROCLOR-1260		25	5		89			3	r 66	۵	34	_		

PROJ_NO: 05483	NSAMPLE	SD-167-0-1			SD-168-0-1			SD-169-0-1			SD-170-0-1			
SDG: 180-28384-1	LAB_ID	180-28384-5			180-28384-6			180-28384-7	-		180-28384-8			
FRACTION: PCB	SAMP_DATE	12/18/2013			12/18/2013			12/18/2013			12/20/2013			
MEDIA: SOIL	QC_TYPE	ΣZ			ΣZ			ΣZ			ΣN			
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG			-
	PCT_SOLIDS 77.6	77.6			78.7			87.6			78.5			
	DUP_OF											•		
PARAMETER		RESULT	ΛαΓ	QLCD	RESULT	VQL	QLCD	RESULT	NOF	arcp	RESULT	VQL	arcd	
AROCLOR-1016		0.4 U	n		0.39	0 e		1 96.0	15 U		0.39	n e		
AROCLOR-1221		0.51	ם		0.5	2 0		0.4	0.45 U		0.51	1 N		
AROCLOR-1232		0.46 U	n		0.45	3 U		4.0	0.41 U		0.45	2 0		-
AROCLOR-1242		0.44	n		0.43	3 U		0.39	n ရ		0.43	3 U		
AROCLOR-1248		0.25 U	D		0.25	10		0.2	0.22 U		0.25	5 U		
AROCLOR-1254		0.38	D.		100	_		2	29		0.38	8 U		
AROCLOR-1260		19			8	-		2	29		2.	2.1 J	Ь	

PROJ_NO: 05483	NSAMPLE	SD-171-0-1			SD-172-0-3			SD-173-0-3			SD-174-0-1			Г
SDG: 180-28384-1	LAB_ID	180-28384-9			180-28384-10	0		180-28384-11	-		180-28384-12	2		
FRACTION: PCB	SAMP_DATE	12/19/2013			12/19/2013			12/19/2013			12/19/2013			
MEDIA: SOIL	QC_TYPE	ΣZ			ΣN			ΝM			ΣZ		,	
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG			
	PCT_SOLIDS 33.2	33.2			53.1			62.0			83.4			
	DUP_OF													
PARAMETER		RESULT	VaL	alcd	RESULT	VQL	QLCD	RESULT	VaL	arcp	RESULT	VQL	QLCD	
AROCLOR-1016		190 [	) U		12	120 U	<u></u>	)5	200 n		0.3	0.37 U		
AROCLOR-1221		240	<u> </u>		15	150 U		79	640 U		0.4	0.48 U		
AROCLOR-1232		210 L	ΩC		13	130 U		35	580 U		0.43	3 U		
AROCLOR-1242		200 U	n		13	130 U		55	250 U		0.4	0.41 U		
AROCLOR-1248		120 U	n c		2	74 U		35	320 U		0.24	74 U		
AROCLOR-1254		180 U	n (		11	110 U		48	480 U	-	150	0		
AROCLOR-1260	,	150000			54000	0		220000	00		350	0		

PROJ_NO: 05483	NSAMPLE	SD-175-0-1			SD-176-0-3			SD-177-0-2			SD-178-0-1		
SDG: 180-28384-1	LAB_ID	180-28384-13	3		180-28384-14	4		180-28384-15	15		180-28384-16		
FRACTION: PCB	SAMP_DATE	12/19/2013			12/19/2013			12/19/2013			12/19/2013		
MEDIA: SOIL	QC_TYPE	ΣZ			ΣN			ΝN			ΣZ		
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG		
	PCT_SOLIDS 70.0	70.0			56.7			28.2			78.6		
	DUP_OF												
PARAMETER		RESULT	Val	arcp	RESULT	NOL	arcp	RESULT	VQL	arcd	RESULT	Val.	arco
AROCLOR-1016		4	4.4 U		0.55	5 U		-	1.1 W	Y	1 66.0		
AROCLOR-1221		5.7	7 U		0.7	7 U			1.4 UJ	٨	0.5		
AROCLOR-1232		5.1	1 0		1 69.0	3 U		1	1.3 UJ	٨	0.45		
AROCLOR-1242		4	4.8 U		9.0	9			1.2 UJ	Y	0.43		
AROCLOR-1248		2.	2.8 U		0.35	5 U		0.0	0.69 UJ	٨	0.25		
AROCLOR-1254		410	0		98	9			28 J	Y	5.3		
AROCLOR-1260		1300	0		88	8			73 J	>	12		

PROJ_NO: 05483	NSAMPLE	SD-179-0-1			SD-180-0-1			SD-181-0-1			SD-182-0-3		
SDG: 180-28384-1	LAB_ID	180-28384-17			180-28384-18	80		180-28384-19	6		180-28384-20		
FRACTION: PCB	SAMP_DATE	12/19/2013			12/19/2013			12/20/2013			12/20/2013		
MEDIA: SOIL	QC_TYPE	ΣN			ΣN			ΣZ			NN		
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG		
	PCT_SOLIDS 69.7	69.7			78.7			57.5			47.9		
	DUP_OF												
PARAMETER		RESULT	VQL	QLCD	RESULT	NOL	arcd	RESULT	VQL	alcd	RESULT	VQL	alcd
AROCLOR-1016		0.44 U	O I		0.79 ר	N 6		210	2100 U		2600	n	
AROCLOR-1221		U 75.0	) )			1 U		2800 ר	n o		3300	n	
AROCLOR-1232		0.51	ם		0.91	1 0		2500	n o		3000	n	
AROCLOR-1242		0.49	D.		0.86	9 N		240	2400 U		2800 U	n	
AROCLOR-1248		0.28	D.		0.5	2 N		140	1400 U		1600 U	n	
AROCLOR-1254		7.6			21	1		210	2100 U	-	2500	n	
AROCLOR-1260		22			95	2		420000	0		780000		

PROJ_NO: 05483	NSAMPLE	SD-183-0-1		:	SD-184-0-2			SD-185-0-1			SD-186-0-1		
SDG: 180-28384-1	LAB_ID	180-28384-21	_		180-28384-22			180-28384-23	<u>г</u>		180-28384-24		
FRACTION: PCB	SAMP_DATE	12/20/2013			12/20/2013			12/20/2013			12/20/2013		
MEDIA: SOIL	QC_TYPE	MN			ΣN			ΣN			ΣN		
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG		
	PCT_SOLIDS 30.1	30.1			67.9			74.6			74.2		
	DUP_OF			:									
PARAMETER		RESULT	Val	arcp	RESULT	VQL	QLCD	RESULT	VQL	OLCD	RESULT	VaL	alcd
AROCLOR-1016		41	410 U		0.91	n		3.0	0.83 U		0.84	n ·	_
AROCLOR-1221	٠	53	530 U		1.2	1.2 U		1	1.1 U		1.1	1.1 U	
AROCLOR-1232		47	470 U			n		3.0	0.96 U		96.0	D.	
AROCLOR-1242		45	450 U		n 66:0	n l		0.91	11 U		0.91	n	
AROCLOR-1248		26	260 U		0.58	n		0.6	0.53 U		0.53	n ı	
AROCLOR-1254		39	390 U		410			18	190		38		
AROCLOR-1260		40000	0		1300			550	0:		09		

PROJ_NO: 05483	NSAMPLE	SD-187-0-1			SD-188-0-1			SD-189-0-1			SD-190-0-1	,		
SDG: 180-28384-1	LAB_ID	180-28384-25	,_		180-28384-26			180-28384-27	7.		180-28384-28	~		
FRACTION: PCB	SAMP_DATE	12/20/2013			12/20/2013			12/20/2013			12/20/2013			
MEDIA: SOIL	QC_TYPE	ΣZ			ΣZ			NN			ΣZ		į	
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG		:	
	PCT_SOLIDS 79.7	79.7		,	88.3			9.06			8.69			
	DUP_OF										:			
PARAMETER		RESULT	VQL	arcd	RESULT	VQL	arcd	RESULT	VQL	arcd	RESULT	VQL	alcd	
AROCLOR-1016		U 87.0	O.		0.5	0.7.U		0.68	n 8		0.88	3 U		
AROCLOR-1221		0.99	D t		6.0	n e		0.87	N 2:		1.:	1.1 U		
AROCLOR-1232		68.0	n		0.81	1 0		0.78	N 8.		,-	l U		
AROCLOR-1242		0.85	D		72.0	7 U		0.75	.e u		1 26.0	، ا ں		
AROCLOR-1248		0.49	n c		0.45	2 U		0.43	3 U		0.56	s U		
AROCLOR-1254		29	-		15	5		1	11		53	~		
AROCLOR-1260		92			49	6		80	82		46	3		

PROJ_NO: 05483	NSAMPLE	SD-191-0-1			SD-192-0-1			SD-193-0-1			SD-194-0-1		,	
SDG: 180-28384-1	LAB_ID	180-28384-29			180-28384-30			180-28384-31	31		180-28384-32	2		
FRACTION: PCB	SAMP_DATE	12/20/2013			12/20/2013			12/20/2013			12/20/2013			
MEDIA: SOIL	QC_TYPE	ΣN			ΝN			ΝN			ΣZ			
	UNITS	UG/KG			UG/KG			UG/KG			UG/KG			
	PCT_SOLIDS 69.9	6.69			76.2			70.8			84.3			
	DUP_OF													
PARAMETER		RESULT	VQL	arcp	RESULT	VQL	arcp	RESULT	VQL	arcd	RESULT	VaL	arcd	
AROCLOR-1016		0.89	ם		0.81	n		0.8	0.88 U		0.73	3 U		
AROCLOR-1221		1.1	<u> </u>			۱ ۱		1	1.1 U		0.94	4 U		
AROCLOR-1232		_	n		0.94	n t			1 0		0.84	74 U		
AROCLOR-1242		0.97	<u></u>		0.89	n e		3.0	0.96 U		0.	0.8 U		
AROCLOR-1248		95.0	n	-	0.52	2 U		0.6	0.56 U		0.47	7 U		
AROCLOR-1254		350	_		13	8		0.8	0.84 U		0.	0.7 U		
AROCLOR-1260		230			20	)			17			16		

# Appendix B

Results as Reported by the Laboratory

Job Number: 180-28384-1

Client Sample ID:

SD-163-0-1

Lab Sample ID:

180-28384-1

Client Matrix:

Sediment

% Moisture:

23.2

Date Sampled: 12/18/2013 1155

Date Received: 12/23/2013 0900

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method: Dilution: 3541 50 Prep Batch:

180-93179

Lab File ID:

D1231007.D

Analysis Date:
Prep Date:

12/31/2013 1456 12/24/2013 0315 Initial Weight/Volume:

30.1 g

Final Weight/Volume: Injection Volume:

0.5 mL 2 uL

			•			
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
Anthracene	udin	97	J	21	220	
Benzo[a]anthracene		530		27	220	
Benzo[b]fluoranthene		700		34	220	
Benzo[k]fluoranthene		260		44	220	
Benzo[g,h,i]perylene		480		22	220	
Benzo[a]pyrene		500		22	220	
Chrysene		710		26	220	
Dibenz(a,h)anthracene		110	J	24	220	
Fluoranthene		1400		23	220	
Fluorene		52	J	29	220	
Indeno[1,2,3-cd]pyrene		390		22	220	
Phenanthrene		990		34	220	
Pyrene		1100		22	220	
Acenaphthene		64	J	21	220	
Acenaphthylene		69	J	25	220	
Naphthalene		ND		19	220	
2-Methylnaphthalene		ND		19	220	
Surrogate		%Rec	Qualifier	Accepta	nce Limits	
Nitrobenzene-d5 (Surr)		64	D	27 - 110	**************************************	
2-Fluorobiphenyl		70	D	28 - 108	1	
Terphenyl-d14 (Surr)		72	D	21 - 130	1	

Job Number: 180-28384-1

Client Sample ID:

SD-164-0-1

Lab Sample ID:

180-28384-2

Client Matrix:

Sediment

% Moisture:

21.2

Date Sampled: 12/18/2013 1210

Date Received: 12/23/2013 0900

8270D LL :	Semivolatile •	Organic Com	pounds by	GC/MS - Low	Level

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

Prep Batch:

Lab File ID:

Dilution:

50

D1231008.D

Analysis Date:

180-93179

Initial Weight/Volume:

30.0 g

12/31/2013 1522

Final Weight/Volume:

0.5 mL

,	
Prep	Date:

12/24/2013 0315

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene	**************************************	200	J	21	210
Benzo[a]anthracene		520		27	210
Benzo[b]fluoranthene		640		33	210
Benzo[k]fluoranthene		420		43	210
Benzo[g,h,i]perylene		760		21	210
Benzo[a]pyrene		660		21	210
Chrysene		760		25	210
Dibenz(a,h)anthracene		130	J	24	210
luoranthene		960		23	210
luorene		52	J	28	210
ndeno[1,2,3-cd]pyrene		560		22	210
henanthrene		370		34	210
⊃yrene		1100		21	210
Acenaphthene		25	J	20	210
Acenaphthylene		150	J	24	210
Naphthalene		34	J	18	210
2-Methylnaphthalene		30	J	19	210
Surrogate		%Rec	Qualifier	Accepta	ance Limits
Nitrobenzene-d5 (Surr)	thatile whitemen: "	74	D	27 - 11	0
2-Fluorobiphenyl		83	D	28 - 10	8
Terphenyl-d14 (Surr)		63	D	21 - 13	0

Client: Tetra Tech, Inc.

Job Number: 180-28384-1

Client Sample ID:

SD-165-0-1

Lab Sample ID:

180-28384-3

Client Matrix:

Sediment

% Moisture:

23.8

Date Sampled: 12/18/2013 1235

Date Received: 12/23/2013 0900

8270D LL Semivolatile	Organia Compoundo	by CC/MC Lavel aval
62/UD LL Semivolatile	Organic Compounds	DV GC/NS - LOW Level

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method: Dilution:

3541

Prep Batch:

180-93179

Lab File ID:

Analysis Date:

50

Initial Weight/Volume:

21 - 130

D1231009.D

Prep Date:

12/31/2013 1549 12/24/2013 0315 Final Weight/Volume: Injection Volume:

30.1 g 0.5 mL

D

2 uL

1-6-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1						
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
Anthracene	ttermetettitetettemmermaansersaanse matematike	110	J	21	220	
Benzo[a]anthracene		830		27	220	
Benzo[b]fluoranthene		1400		34	220	
Benzo[k]fluoranthene		590		44	220	
Benzo[g,h,i]perylene		890		22	220	
Benzo[a]pyrene		920		22	220	
Chrysene		1500		26	220	
Dibenz(a,h)anthracene		200	J	24	220	
Fluoranthene		3500		23	220	
Fluorene		55	J	29	220	
Indeno[1,2,3-cd]pyrene		790		22	220	
Phenanthrene		2800		35	220	
Pyrene		2500		22	220	
Acenaphthene		47	J	21	220	
Acenaphthylene		36	J	25	220	
Naphthalene		95	J	19	220	
2-Methylnaphthalene		25	J	20	220	
Surrogate		%Rec	Qualifier	Accepta	ance Limits	
Nitrobenzene-d5 (Surr)	**************************************	81	D	27 - 110	0	
2-Fluorobiphenyl		80	D	28 - 108	3	

67

Terphenyl-d14 (Surr)

Client: Tetra Tech, Inc.

Job Number: 180-28384-1

Client Sample ID:

SD-166-0-1

Lab Sample ID:

180-28384-4

Client Matrix:

Sediment

% Moisture:

20.0

Date Sampled: 12/18/2013 1305

Date Received: 12/23/2013 0900

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method: Dilution:

3541 100

Prep Batch:

180-93179

Lab File ID:

D1231010.D

ΧD

ΧD

Initial Weight/Volume:

30.0 g

Analysis Date: Prep Date:

2-Fluorobiphenyl

Terphenyl-d14 (Surr)

12/31/2013 1615 12/24/2013 0315 Final Weight/Volume: Injection Volume:

28 - 108

21 - 130

1.0 mL 2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene	- ** *********************************	ND		82	840
Benzo[a]anthracene		130	J	100	840
Benzo[b]fluoranthene		ND		130	840
Benzo[k]fluoranthene		ND		170	840
Benzo[g,h,i]perylene		130	J	83	840
Benzo[a]pyrene		120	J	83	840
Chrysene		180	J	99	840
Dibenz(a,h)anthracene		100	J	93	840
Fluoranthene		180	J	89	840
Fluorene		ND		110	840
Indeno[1,2,3-cd]pyrene		ND		86	840
Phenanthrene		ND		130	840
Pyrene		210	J	84	840
Acenaphthene		ND		80	840
Acenaphthylene		ND		96	840
Naphthalene		ND		72	840
2-Methylnaphthalene		ND		75	840
Surrogate		%Rec	Qualifier	Accepta	ance Limits
Nitrobenzene-d5 (Surr)		0	ΧD	27 - 110	)

0

0

Client: Tetra Tech, Inc.

Job Number: 180-28384-1

Client Sample ID:

SD-167-0-1

Lab Sample ID:

180-28384-5

Client Matrix:

Sediment

% Moisture:

22.4

Date Sampled: 12/18/2013 1510

Date Received: 12/23/2013 0900

8270D	11	Semivolatile	Organic	Compounds	hν	GC/MS - Low Level	

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

Prep Batch:

180-93179

Lab File ID:

Dilution:

100

D1231011.D

Initial Weight/Volume:

30.0 g

Analysis Date:

12/31/2013 1641

Final Weight/Volume:

0.5 mL

Prep Date:

12/24/2013 0315

Injection Volume:

010 0010		Hijeo	don volume.	Z UL
DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
erromanierromanierromanierromanierromanierromanierromanierromanierromanierromanierromanierromanierromanierroma	44	J	42	430
	270	J	54	430
	430		68	430
	170	J	87	430
	340	J	43	430
	280	J	43	430
	440		51	430
	ND		48	430
	800		46	430
	ND		57	430
	300	J	44	430
	340	J	68	430
	590		43	430
	ND		41	430
	ND		49	430
	ND		37	430
	ND		39	430
	%Rec	Qualifier	Accepta	nce Limits
econoccione economica — no - emilio existe i consistente consistente econoccione econoccio	0	ΧD	27 - 110	· · · · · · · · · · · · · · · · · · ·
	0	ΧD	28 - 108	
	0	X D	21 - 130	
		DryWt Corrected: Y  Result (ug/Kg)  44  270  430  170  340  280  440  ND  800  ND  300  340  590  ND  ND  ND  ND  ND  ND  ND  ND  ND  N	DryWt Corrected: Y  Result (ug/Kg)  Qualifier  44  J  270  J  430  170  J  340  J  280  J  440  ND  800  ND  300  J  340  J  590  ND  ND  ND  ND  ND  ND  ND  ND  ND  N	DryWt Corrected: Y Result (ug/Kg) Qualifier MDL  44 J 42 270 J 54 430 68 170 J 87 340 J 43 280 J 43 440 51 ND 48 800 46 ND 57 300 J 44 340 J 68 590 43 ND J 68 590 43 ND 41 ND 49 ND 37 ND 39  %Rec Qualifier Accepta 0 X D 27 - 110 0 X D 28 - 108

Client: Tetra Tech, Inc. Job Number: 180-28384-1

Client Sample ID:

SD-168-0-1

Lab Sample ID: Client Matrix:

180-28384-6

Sediment

% Moisture:

21.3

Date Sampled: 12/18/2013 1545

Date Received: 12/23/2013 0900

8270D LL Semivolatile	Organic Compounds	by CC/MS - Low Lovel
62/UD LL Semivolatile	Organic Compounds	DV GC/IVIS - LOW Level

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541 100

Prep Batch:

180-93179

Lab File ID:

Dilution:

Initial Weight/Volume:

D1231012.D

Analysis Date:

12/31/2013 1708

Final Weight/Volume:

30.1 g 0.5 mL

Prep Date:

12/24/2013 0315

Injection Volume:

•			•		
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene	Print ten assur Luntesstilletontelari vindermiselini arritessatess	180	J	41	420
Benzo[a]anthracene		470		53	420
Benzo[b]fluoranthene		250	J	66	420
Benzo[k]fluoranthene		530		85	420
Benzo[g,h,i]perylene		410	J	42	420
Benzo[a]pyrene		390	J	42	420
Chrysene		530		50	420
Dibenz(a,h)anthracene		88	J	47	420
Fluoranthene		1000		45	420
Fluorene		86	J	56	420
Indeno[1,2,3-cd]pyrene		370	J	44	420
Phenanthrene		710		67	420
Pyrene		830		43	420
Acenaphthene		110	J	41	420
Acenaphthylene		ND		48	420
Naphthalene		36	J	36	420
2-Methylnaphthalene		ND		38	420
Surrogate		%Rec	Qualifier	Acceptar	nce Limits
Nitrobenzene-d5 (Surr)	**************************************	0	X D	27 - 110	gggygger in interested and the second and interested and in in in in-
2-Fluorobiphenyl		0	ΧD	28 - 108	
Terphenyl-d14 (Surr)		0	ΧD	21 - 130	

Job Number: 180-28384-1

Client Sample ID:

SD-169-0-1

Lab Sample ID:

180-28384-7

Client Matrix:

Sediment

% Moisture:

12.4

Date Sampled: 12/18/2013 1610

Date Received: 12/23/2013 0900

8270D LL Semivolatile	Organic Compounds b	v GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method: Dilution:

3541

Prep Batch:

180-93179

Lab File ID:

50

Initial Weight/Volume:

D1231013.D

Analysis Date:

12/31/2013 1734

Final Weight/Volume:

30.1 g 0.5 mL

Analysis Date	;
Prep Date:	

12/24/2013 0315

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene	**************************************	ND	modifiabrenia - modifiable della sociazioni e minoriare si e i i i i i	19	190
Benzo[a]anthracene		76	J	24	190
Benzo[b]fluoranthene		120	J *	30	190
Benzo[k]fluoranthene		43	J *	38	190
Benzo[g,h,i]perylene		81	J *	19	190
Benzo[a]pyrene		69	J *	19	190
Chrysene		120	J	23	190
Dibenz(a,h)anthracene		ND	*	21	190
Fluoranthene		180	J	20	190
Fluorene		ND		25	190
Indeno[1,2,3-cd]pyrene		66	J *	20	190
Phenanthrene		47	J	30	190
Pyrene		160	J	19	190
Acenaphthene		30	J	18	190
Acenaphthylene		ND		22	190
Naphthalene		ND		16	190
2-Methylnaphthalene		ND		17	190
Surrogate		%Rec	Qualifier	Acceptar	nce Limits
Nitrobenzene-d5 (Surr)		68	D	27 - 110	gi (
2-Fluorobiphenyl		51	D	28 - 108	
Terphenyl-d14 (Surr)		55	D	21 - 130	

Client: Tetra Tech, Inc. Job Number: 180-28384-1

Client Sample ID:

SD-170-0-1

Lab Sample ID:

180-28384-8

Client Matrix:

Sediment

% Moisture:

21.5

Date Sampled: 12/20/2013 1515

Date Received: 12/23/2013 0900

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

Lab File ID:

Prep Batch:

180-93179

D1231014.D

Dilution:

50

Initial Weight/Volume: Final Weight/Volume:

30.0 g

Analysis Date:

12/31/2013 1800

0.5 mL

Prep Date:

12/24/2013 0315

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		110	J	21	210
Benzo[a]anthracene		320		27	210
Benzo[b]fluoranthene		320	*	33	210
Benzo[k]fluoranthene		78	J *	43	210
Benzo[g,h,i]perylene		170	J *	21	210
Benzo[a]pyrene		250	*	21	210
Chrysene		280		25	210
Dibenz(a,h)anthracene		32	J *	24	210
Fluoranthene		630		23	210
Fluorene		ND		28	210
Indeno[1,2,3-cd]pyrene		150	J*	22	210
Phenanthrene		300		34	210
Pyrene		480		21	210
Acenaphthene		ND		20	210
Acenaphthylene		58	J	24	210
Naphthalene		ND		18	210
2-Methylnaphthalene		ND		19	210
Surrogate		%Rec	Qualifier	Accepta	nce Limits
Nitrobenzene-d5 (Surr)		65	D	27 - 110	**************************************
2-Fluorobiphenyl		67	D	28 - 108	
Terphenyl-d14 (Surr)		57	D	21 - 130	

Job Number: 180-28384-1

Client Sample ID:

SD-171-0-1

Lab Sample ID:

180-28384-9

Client Matrix:

Sediment

% Moisture:

66.8

Date Sampled: 12/19/2013 1045

Date Received: 12/23/2013 0900

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

Prep Batch:

Lab File ID:

Dilution:

180-93179

D1231015.D

5.0

Initial Weight/Volume: Final Weight/Volume:

30.0 g

Analysis Date:

12/31/2013 1826

0.5 mL

Prep Date:

12/24/2013 0315

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
Anthracene		160		4.9	50	**********
Benzo[a]anthracene		620		6.3	50	
Benzo[b]fluoranthene		720	*	7.9	50	
Benzo[k]fluoranthene		250	*	10	50	
Benzo[g,h,i]perylene		630	*	5.0	50	
Benzo[a]pyrene		680	*	5.0	50	
Chrysene		720		6.0	50	
Dibenz(a,h)anthracene		150	*	5.6	50	
Fluoranthene		910		5.4	50	
Fluorene		35	J	6.6	50	
Indeno[1,2,3-cd]pyrene		530	*	5.2	50	
Phenanthrene		430		8.0	50	
Pyrene		1100		5.1	50	
Acenaphthene		28	J	4.8	50	
Acenaphthylene		71		5.8	50	
Naphthalene		24	J	4.3	50	
2-Methylnaphthalene		14	J	4.5	50	
Surrogate		%Rec	Qualifier	Accept	ance Limits	
Nitrobenzene-d5 (Surr)		54	**************************************	27 - 11	0	0000000000
2-Fluorobiphenyl		58		28 - 10	8	
Terphenyl-d14 (Surr)		53		21 - 13	0	

Job Number: 180-28384-1

Client Sample ID:

SD-172-0-3

Lab Sample ID:

180-28384-10

Client Matrix:

Sediment

% Moisture:

46.9

Date Sampled: 12/19/2013 1110

Date Received: 12/23/2013 0900

8270D LL Semivolatile Orga	nic Compounds !	by GC/MS -	Low Level
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Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

180-93179

Lab File ID:

Dilution:

Prep Batch:

D1231016.D

50

Initial Weight/Volume:

Final Weight/Volume:

30.1 g

Analysis Date: Prep Date:

12/31/2013 1853 12/24/2013 0315

Injection Volume:

0.5 mL 2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		1100		31	310
Benzo[a]anthracene		4600		39	310
Benzo[b]fluoranthene		5300	*	49	310
Benzo[k]fluoranthene		1800	*	63	310
Benzo[g,h,i]perylene		3300	*	31	310
Benzo[a]pyrene		4200	*	31	310
Chrysene		5000		37	310
Dibenz(a,h)anthracene		870	*	35	310
Fluoranthene		9600		33	310
Fluorene		460		41	310
ndeno[1,2,3-cd]pyrene		3100	*	32	310
Phenanthrene		5000		50	310
Pyrene		7400		32	310
Acenaphthene		550		30	310
Acenaphthylene		160	J	36	310
Naphthalene		89	J	27	310
2-Methylnaphthalene		50	J	28	310

Surrogate	%Rec	Qualifier	Acceptance Limits
Nitrobenzene-d5 (Surr)	54	D	27 - 110
2-Fluorobiphenyl	53	D	28 - 108
Terphenyl-d14 (Surr)	47	D	21 - 130

Job Number: 180-28384-1

Client Sample ID:

SD-173-0-3

Lab Sample ID:

180-28384-11

Client Matrix:

Sediment

% Moisture:

38.0

Date Sampled: 12/19/2013 1130

Date Received: 12/23/2013 0900

#### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

Prep Batch:

Lab File ID:

D1231017.D

Dilution:

180-93179

50

Initial Weight/Volume:

30.0 g

Analysis Date:

12/31/2013 1919

Final Weight/Volume:

0.5 mL

Prep Date:

12/24/2013 0315

Injection Volume:

Analyte Anthracene Benzo[a]anthracene Benzo[b]fluoranthene Benzo[k]fluoranthene	DryWt Corrected: Y	Result (ug/Kg) 580 1700	Qualifier	MDL 26	RL
Benzo[a]anthracene Benzo[b]fluoranthene				~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	**************************************
Benzo[b]fluoranthene		1700		20	270
				34	270
Ponzolkifluoranthono		2100	*	42	270
Denzo[k]nuorannene		630	*	54	270
Benzo[g,h,i]perylene		1800	*	27	270
Benzo[a]pyrene		1900	*	27	270
Chrysene		2000		32	270
Dibenz(a,h)anthracene		390	*	30	270
Fluoranthene		3000		29	270
Fluorene		160	J	35	270
Indeno[1,2,3-cd]pyrene		1500	*	28	270
Phenanthrene		1100		43	270
Pyrene		3400		27	270
Acenaphthene		150	J	26	270
Acenaphthylene		430		31	270
Naphthalene		94	J	23	270
2-Methylnaphthalene		55	J	24	270
Surrogate		%Rec	Qualifier	Accepta	nce Limits
Nitrobenzene-d5 (Surr)		65	D	27 - 110	)
2-Fluorobiphenyl		67	D	28 - 108	}
Terphenyl-d14 (Surr)		59	D	21 - 130	)

Client Sample ID:

SD-174-0-1

Lab Sample ID:

180-28384-12

Client Matrix:

Sediment

% Moisture:

16.6

Date Sampled: 12/19/2013 1340

Date Received: 12/23/2013 0900

#### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

Prep Batch:

180-93179

Lab File ID:

D1231018.D

Dilution:

50

Initial Weight/Volume: Final Weight/Volume:

30.0 g

Analysis Date: Prep Date:

12/31/2013 1946

12/24/2013 0315

Injection Volume:

0.5 mL 2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		610	······	20	200
Benzo[a]anthracene		2200		25	200
Benzo[b]fluoranthene		2500	*	31	200
Benzo[k]fluoranthene		780	*	40	200
Benzo[g,h,i]perylene		1300	*	20	200
Benzo[a]pyrene		1900	*	20	200
Chrysene		2400		24	200
Dibenz(a,h)anthracene		360	*	22	200
Fluoranthene		5800		21	200
-luorene		250		26	200
ndeno[1,2,3-cd]pyrene		1300	*	21	200
Phenanthrene		3800		32	200
Pyrene		4000		20	200
Acenaphthene		230		19	200
Acenaphthylene		140	J	23	200
Naphthalene		160	J	17	200
2-Methylnaphthalene		47	J	18	200
Surrogate		%Rec	Qualifier	Acceptar	nce Limits
Nitrobenzene-d5 (Surr)		65	D	27 - 110	
?-Fluorobiphenyl		69	D	28 - 108	
Terphenyl-d14 (Surr)		58	D	21 - 130	

Client Sample ID:

SD-175-0-1

Lab Sample ID:

180-28384-13

Client Matrix:

Sediment

% Moisture:

30.0

Date Sampled: 12/19/2013 1430

Date Received: 12/23/2013 0900

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

Lab File ID:

D1231019.D

Dilution:

Prep Batch:

180-93179

Initial Weight/Volume:

30.0 g

Analysis Date:

50

12/31/2013 2012

Final Weight/Volume:

0.5 mL

Prep Date:

12/24/2013 0315

Injection Volume:

riep Date. 12/24	W2013 0313		injection volume.			
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
Anthracene		2100		23	240	***************************************
Benzo[a]anthracene		5700		30	240	
Benzo[b]fluoranthene		6000	*	37	240	
Benzo[k]fluoranthene		2500	*	48	240	
Benzo[g,h,i]perylene		3700	*	24	240	
Benzo[a]pyrene		5400	*	24	240	
Chrysene		6100		28	240	
Dibenz(a,h)anthracene		1100	*	26	240	
Fluoranthene		13000		25	240	
Fluorene		1000		31	240	
Indeno[1,2,3-cd]pyrene		3600	*	25	240	
Phenanthrene		9000		38	240	
Pyrene		10000		24	240	
Acenaphthene		1500		23	240	
Acenaphthylene		310		27	240	
Naphthalene		480		21	240	
2-Methylnaphthalene		210	J	21	240	
Surrogate		%Rec	Qualifier	Accepta	ınce Limits	
Nitrobenzene-d5 (Surr)		63	D	27 - 110	)	***************************************
2-Fluorobiphenyl		62	D	28 - 108	3	
Terphenyl-d14 (Surr)		64	D	21 - 130	)	

Job Number: 180-28384-1

Client Sample ID:

SD-176-0-3

Lab Sample ID:

180-28384-14

Client Matrix:

Sediment

% Moisture:

43.3

Date Sampled: 12/19/2013 1505

Date Received: 12/23/2013 0900

#### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

180-93179

Lab File ID:

D1231020.D

Prep Batch:

Dilution:

50

Initial Weight/Volume:

30.1 g

Analysis Date:

12/31/2013 2038

Final Weight/Volume:

0.5 mL

Prep Date:

12/24/2013 0315

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		390		29	290
Benzo[a]anthracene		1600		37	290
Benzo[b]fluoranthene		2700	*	46	290
Benzo[k]fluoranthene		710	*	59	290
Benzo[g,h,i]perylene		1500	*	29	290
Benzo[a]pyrene		1800	*	29	290
Chrysene		2200		35	290
Dibenz(a,h)anthracene		320	*	33	290
Fluoranthene		4200		31	290
Fluorene		160	J	39	290
ndeno[1,2,3-cd]pyrene		1400	*	30	290
Phenanthrene		1700		47	290
Pyrene		3000		30	290
Acenaphthene		130	J	28	290
Acenaphthylene		200	J	34	290
Naphthalene		51	J	25	290
2-Methylnaphthalene		30	J	26	290
Surrogate		%Rec	Qualifier	Accepta	nce Limits
Nitrobenzene-d5 (Surr)		57	D	27 - 110	
2-Fluorobiphenyl		57	D	28 - 108	
Terphenyl-d14 (Surr)		45	D	21 - 130	

Job Number: 180-28384-1

Client Sample ID:

SD-177-0-2

Lab Sample ID:

180-28384-15

Client Matrix:

Sediment

% Moisture:

71.8

Date Sampled: 12/19/2013 1540

Date Received: 12/23/2013 0900

### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

Prep Batch:

Lab File ID:

D1231021.D

Dilution:

180-93179

50

Initial Weight/Volume:

30.2 g

Analysis Date:

Final Weight/Volume:

0.5 mL

Prep Date:

12/31/2013 2104 12/24/2013 0315

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		3300	***************************************	57	590
Benzo[a]anthracene		18000		74	590
Benzo[b]fluoranthene		27000	*	92	590
Benzo[k]fluoranthene		10000	*	120	590
Benzo[g,h,i]perylene		15000	*	58	590
Benzo[a]pyrene		21000	*	59	590
Chrysene		26000		70	590
Dibenz(a,h)anthracene		3500	*	65	590
Fluoranthene		68000		63	590
Fluorene		2200		77	590
Indeno[1,2,3-cd]pyrene		14000	*	60	590
Phenanthrene		39000		93	590
Pyrene		45000		59	590
Acenaphthene		2200		56	590
Acenaphthylene		510	J	67	590
Naphthalene		120	J	51	590
2-Methylnaphthalene		300	J	53	590
Surrogate		%Rec	Qualifier	Accepta	nce Limits
Nitrobenzene-d5 (Surr)		40	D	27 - 110	
2-Fluorobiphenyl		42	D	28 - 108	
Terphenyl-d14 (Surr)		39	D	21 - 130	

Client Sample ID:

SD-178-0-1

Lab Sample ID:

180-28384-16

Client Matrix:

Sediment

% Moisture:

21.4

Date Sampled: 12/19/2013 1600

Date Received: 12/23/2013 0900

8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Prep Method:

3541

Prep Batch:

Lab File ID:

D1231022.D

Dilution:

180-93179

Initial Weight/Volume:

30.0 g

50

Final Weight/Volume:

0.5 mL

Analysis Date: Prep Date:

12/31/2013 2131 12/24/2013 0315

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		30	J	21	210
Benzo[a]anthracene		140	J	27	210
Benzo[b]fluoranthene		200	J *	33	210
Benzo[k]fluoranthene		86	J *	43	210
Benzo[g,h,i]perylene		140	J *	21	210
Benzo[a]pyrene		180	J *	21	210
Chrysene		190	J	25	210
Dibenz(a,h)anthracene		32	J *	24	210
Fluoranthene		260		23	210
Fluorene		ND		28	210
Indeno[1,2,3-cd]pyrene		110	J *	22	210
Phenanthrene		81	J	34	210
Pyrene		210		21	210
Acenaphthene		ND		20	210
Acenaphthylene		48	J	24	210
Naphthalene		ND		18	210
2-Methylnaphthalene		ND		19	210

Client Sample ID:

SD-179-0-1

Lab Sample ID:

180-28384-17

Client Matrix:

Sediment

% Moisture:

30.3

Date Sampled: 12/19/2013 1645

Date Received: 12/23/2013 0900

Analysis Method:

8270D LL

Analysis Batch:

180-93613

Instrument ID:

732

Acceptance Limits

27 - 110 28 - 108

21 - 130

Prep Method:

3541

Prep Batch:

Lab File ID:

D1231023.D

Dilution:

180-93179

Initial Weight/Volume:

30.0 g

Analysis Date:

2.0 12/31/2013 2157

Final Weight/Volume:

0.5 mL

Prep Date:

Surrogate

Nitrobenzene-d5 (Surr)

Terphenyl-d14 (Surr)

2-Fluorobiphenyl

12/24/2013 0353

Injection Volume:

2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		41		0.94	9.6
Benzo[a]anthracene		120		1.2	9.6
Benzo[b]fluoranthene		290	*	1.5	9.6
Benzo[k]fluoranthene		120	*	1.9	9.6
Benzo[g,h,i]perylene		150	*	0.95	9.6
Benzo[a]pyrene		200	*	0.96	9.6
Chrysene		230		1.1	9.6
Dibenz(a,h)anthracene		32	*	1.1	9.6
Fluoranthene		410		1.0	9.6
Fluorene		13		1.3	9.6
ndeno[1,2,3-cd]pyrene		150	*	0.99	9.6
Phenanthrene		150		1.5	9.6
<sup>o</sup> yrene		290		0.97	9.6
Acenaphthene		18		0.92	9.6
Acenaphthylene		14		1.1	9.6
Naphthalene		4.7	J	0.82	9.6
2-Methylnaphthalene		2.5	J	0.86	9.6

Qualifier

%Rec

63

61

46

Client Sample ID:

SD-180-0-1

Lab Sample ID:

180-28384-18

Client Matrix:

Sediment

% Moisture:

21.3

Date Sampled: 12/19/2013 1625

Date Received: 12/23/2013 0900

8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93710

Instrument ID:

732

Prep Method:

3541

Lab File ID:

Prep Batch:

180-93179

D0102001.D

Dilution:

100

Initial Weight/Volume:

30.1 g

Analysis Date:

Final Weight/Volume:

0.5 mL

Prep Date:

01/02/2014 1905 12/24/2013 0353

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		390	J	41	420
Benzo[a]anthracene		2600		53	420
Benzo[b]fluoranthene		4300		66	420
Benzo[k]fluoranthene		1900		85	420
Benzo[g,h,i]perylene		2600		42	420
Benzo[a]pyrene		2900		42	420
Chrysene		3700		50	420
Dibenz(a,h)anthracene		490		47	420
Fluoranthene		6200		45	420
Fluorene		110	J	56	420
Indeno[1,2,3-cd]pyrene		2200		44	420
Phenanthrene		2600		67	420
Pyrene		5900		43	420
Acenaphthene		70	J	41	420
Acenaphthylene		76	J	48	420
Naphthalene		ND		36	420
2-Methylnaphthalene		ND		38	420
Surrogate		%Rec	Qualifier	Accepta	nce Limits
Nitrobenzene-d5 (Surr)		0	ΧD	27 - 110	
2-Fluorobiphenyl		0	ΧD	28 - 108	
Terphenyl-d14 (Surr)		0	ΧD	21 - 130	

Client Sample ID:

SD-181-0-1

Lab Sample ID:

180-28384-19

Client Matrix:

Sediment

% Moisture:

42.5

Date Sampled: 12/20/2013 1000

Date Received: 12/23/2013 0900

8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93710

Instrument ID:

732

Prep Method:

3541

Prep Batch:

Lab File ID:

D0102002.D

Dilution:

180-93179

Initial Weight/Volume:

Injection Volume:

21 - 130

30.0 g

Analysis Date:

100

Final Weight/Volume:

0.5 mL 2 uL

Prep Date:

01/02/2014 1932 12/24/2013 0353

	2010 0000	•					
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL		
Anthracene		7100		57	580		
Benzo[a]anthracene		20000		73	580		
Benzo[b]fluoranthene		21000		91	580		
Benzo[k]fluoranthene		8800		120	580		
Benzo[g,h,i]perylene		14000		58	580		
Benzo[a]pyrene		17000		58	580		
Chrysene		20000		69	580		
Dibenz(a,h)anthracene		3100		64	580		
Fluoranthene		35000		62	580		
Fluorene		4100		76	580		
Indeno[1,2,3-cd]pyrene		12000		60	580		
Phenanthrene		28000		92	580		
Pyrene		37000		59	580		
Acenaphthene		3900		56	580		
Acenaphthylene		350	J	66	580		
Naphthalene		640		50	580		
2-Methylnaphthalene		440	J	52	580		
Surrogate		%Rec	Qualifier	Acceptance Limits			
Nitrobenzene-d5 (Surr)		0	ΧD	27 - 110			
2-Fluorobiphenyl		0	ΧD	28 - 108			
·							

ΧD

Terphenyl-d14 (Surr)

Job Number: 180-28384-1

Client Sample ID:

SD-182-0-3

Lab Sample ID:

180-28384-20

Client Matrix:

Sediment

% Moisture:

52.1

Date Sampled: 12/20/2013 1020

Date Received: 12/23/2013 0900

# 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93710

Instrument ID:

732

Prep Method:

3541

Lab File ID:

Prep Batch:

180-93179

D0102003.D

Dilution:

100

Initial Weight/Volume:

30.1 g

Analysis Date:

Final Weight/Volume:

21 - 130

0.5 mL

Prep Date:

01/02/2014 1958 12/24/2013 0353

Injection Volume:

2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		3000		68	700
Benzo[a]anthracene		10000		87	700
Benzo[b]fluoranthene		13000		110	700
Benzo[k]fluoranthene		3500		140	700
Benzo[g,h,i]perylene		9200		69	700
Benzo[a]pyrene		9500		69	700
Chrysene		11000		83	700
Dibenz(a,h)anthracene		2200		77	700
Fluoranthene		18000		74	700
Fluorene		1200		91	700
Indeno[1,2,3-cd]pyrene		7300		72	700
Phenanthrene		12000		110	700
Pyrene		18000		70	700
Acenaphthene		1600		67	700
Acenaphthylene		230	J	80	700
Naphthalene		450	J	60	700
2-Methylnaphthalene		240	J	62	700
Surrogate		%Rec	Qualifier	Acceptance Limits	
Nitrobenzene-d5 (Surr)		0	ΧD	27 - 110	······································
2-Fluorobiphenyl		0	ΧD	28 - 108	

0

ΧD

Terphenyl-d14 (Surr)

Job Number: 180-28384-1

Client Sample ID:

SD-183-0-1

Lab Sample ID:

180-28384-21

Client Matrix:

Sediment

% Moisture:

69.9

Date Sampled: 12/20/2013 1040

Date Received: 12/23/2013 0900

#### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

Prep Batch:

180-93531

Lab File ID:

S0101005.D

Dilution:

20

Initial Weight/Volume:

30.1 g

Analysis Date: Prep Date:

01/01/2014 0915 12/31/2013 0340

Final Weight/Volume: Injection Volume:

0.5 mL 2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene	/// // · · · · · · · · · · · · · · · ·	5600	***************************************	22	220
Benzo[a]anthracene		15000		28	220
Benzo[b]fluoranthene		15000		35	220
Benzo[k]fluoranthene		6200		45	220
Benzo[g,h,i]perylene		11000		22	220
Benzo[a]pyrene		13000		22	220
Chrysene		16000		26	220
Dibenz(a,h)anthracene		3100		25	220
Fluoranthene		33000		24	220
Fluorene		2400		29	220
Indeno[1,2,3-cd]pyrene		9700		23	220
Phenanthrene		17000		35	220
Pyrene		19000		22	220
Acenaphthene		3000		21	220
Acenaphthylene		320		25	220
Naphthalene		480		19	220
2-Methylnaphthalene		280		20	220

Surrogate	%Rec	Qualifier	Acceptance Limits
Nitrobenzene-d5 (Surr)	51	D	27 - 110
2-Fluorobiphenyl	61	D	28 - 108
Terphenyl-d14 (Surr)	54	D	21 - 130

Client Sample ID:

SD-184-0-2

Lab Sample ID:

180-28384-22

Client Matrix:

Sediment

% Moisture:

32.1

Date Sampled: 12/20/2013 1130

Date Received: 12/23/2013 0900

8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

Prep Batch:

180-93531

Lab File ID:

S0101006.D

Dilution:

20

Initial Weight/Volume:

Analysis Date:

Final Weight/Volume:

30.1 g 0.5 mL

Prep Date:

01/01/2014 0942 12/31/2013 0340

Injection Volume:

2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		2200	(consequences)	9.6	98
Benzo[a]anthracene		4500		12	98
Benzo[b]fluoranthene		4700		15	98
Benzo[k]fluoranthene		1700		20	98
Benzo[g,h,i]perylene		3400		9.7	98
Benzo[a]pyrene		3800		9.8	98
Chrysene		4900		12	98
Dibenz(a,h)anthracene		930		11	98
Fluoranthene		11000		10	98
Fluorene		1100		13	98
Indeno[1,2,3-cd]pyrene		3000		10	98
Phenanthrene		6400		16	98
Pyrene		6100		9.9	98
Acenaphthene		920		9.4	98
Acenaphthylene		560		11	98
Naphthalene		340		8.4	98
2-Methylnaphthalene		170		8.8	98
Surrogate		%Rec	Qualifier	Acceptano	ce Limits
Nitrobenzene-d5 (Surr)		62	D	27 - 110	(A) (A) (A) (A) (A) (A) (A) (A) (A) (A)
2-Fluorobiphenyl		70	D	28 - 108	
Terphenyl-d14 (Surr)		61	D	21 - 130	

Job Number: 180-28384-1

Client Sample ID:

SD-185-0-1

Lab Sample ID:

180-28384-23

Client Matrix:

Sediment

% Moisture:

25.4

Date Sampled: 12/20/2013 1200

Date Received: 12/23/2013 0900

### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

Qualifier

D

D

D

71

Prep Method:

3541

Lab File ID:

S0101007.D

Dilution:

Surrogate

Nitrobenzene-d5 (Surr)

Terphenyl-d14 (Surr)

2-Fluorobiphenyl

20

Prep Batch:

180-93531

Initial Weight/Volume:

30.0 g

Analysis Date: Prep Date:

01/01/2014 1008 12/31/2013 0340 Final Weight/Volume: Injection Volume:

1.0 mL 2 uL

Acceptance Limits

27 - 110

28 - 108

21 - 130

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
Anthracene		230	······································	18	180	***************************************
Benzo[a]anthracene		1300		22	180	
Benzo[b]fluoranthene		1500		28	180	
Benzo[k]fluoranthene		570		36	180	
Benzo[g,h,i]perylene		1100		18	180	
Benzo[a]pyrene		1200		18	180	
Chrysene		1600		21	180	
Dibenz(a,h)anthracene		270		20	180	
Fluoranthene		3600		19	180	
Fluorene		87	J	24	180	
Indeno[1,2,3-cd]pyrene		1000		18	180	
Phenanthrene		1100		28	180	
Pyrene		1900		18	180	
Acenaphthene		53	J	17	180	
Acenaphthylene		94	J	20	180	
Naphthalene		43	J	15	180	
2-Methylnaphthalene		21	J	16	180	

%Rec

62

74

61

Job Number: 180-28384-1

Client Sample ID:

SD-186-0-1

Lab Sample ID:

180-28384-24

Client Matrix:

Sediment

% Moisture:

25.8

Date Sampled: 12/20/2013 1220

Date Received: 12/23/2013 0900

### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

Prep Batch:

180-93531

Lab File ID:

S0101008.D

Dilution:

20

Initial Weight/Volume: Final Weight/Volume:

30.0 g

Analysis Date: Prep Date:

01/01/2014 1035 12/31/2013 0340

Injection Volume:

2.0 mL 2 uL

1201/2010 00 10			,			
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
Anthracene		96	J	35	360	***************************************
Benzo[a]anthracene		280	J	45	360	
Benzo[b]fluoranthene		540		57	360	
Benzo[k]fluoranthene		150	J	73	360	
Benzo[g,h,i]perylene		350	J	36	360	
Benzo[a]pyrene		290	J	36	360	
Chrysene		520		43	360	
Dibenz(a,h)anthracene		67	J	40	360	
Fluoranthene		1100		38	360	
Fluorene		ND		47	360	
Indeno[1,2,3-cd]pyrene		300	J	37	360	
Phenanthrene		400		57	360	
Pyrene		590		36	360	
Acenaphthene		ND		35	360	
Acenaphthylene		50	J	41	360	
Naphthalene		ND		31	360	
2-Methylnaphthalene		ND		32	360	
Surrogate		%Rec	Qualifier	Accepta	ance Limits	
Nitrobenzene-d5 (Surr)	**************************************	76	D	27 - 110	)	1000.00.00.00.00.00.00.00.00.00.00.00.00
2-Fluorobiphenyl (		90	D	28 - 108	3	
Terphenyl-d14 (Surr)		76	D	21 - 130	)	
, ,						

Job Number: 180-28384-1

Client Sample ID:

SD-187-0-1

Lab Sample ID:

180-28384-25

Client Matrix:

Sediment

% Moisture:

20.3

Date Sampled: 12/20/2013 1245

Date Received: 12/23/2013 0900

8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

Lab File ID:

S0101009.D

Dilution:

Prep Batch:

180-93531

Initial Weight/Volume:

20

01/01/2014 1101

Final Weight/Volume:

28 - 108

21 - 130

30.1 g 3.0 mL

Analysis Date: Prep Date:

2-Fluorobiphenyl

Terphenyl-d14 (Surr)

12/31/2013 0340

Injection Volume:

2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		260	J	49	500
Benzo[a]anthracene		1100		63	500
Benzo[b]fluoranthene		1800		79	500
Benzo[k]fluoranthene		720		100	500
Benzo[g,h,i]perylene		1600		50	500
Benzo[a]pyrene		1400		50	500
Chrysene		1700		60	500
Dibenz(a,h)anthracene		340	J	56	500
Fluoranthene		2900		54	500
Fluorene		82	J	66	500
Indeno[1,2,3-cd]pyrene		1300		52	500
Phenanthrene		850		80	500
Pyrene		1900		51	500
Acenaphthene		ND		48	500
Acenaphthylene		190	J	57	500
Naphthalene		ND		43	500
2-Methylnaphthalene		48	J .	45	500
Surrogate		%Rec	Qualifier	Acceptance Limits	
Nitrobenzene-d5 (Surr)		77	D	27 - 110	

D

D

92

83

Job Number: 180-28384-1

Client Sample ID:

SD-188-0-1

Lab Sample ID:

180-28384-26

Client Matrix:

Sediment

% Moisture:

11.7

Date Sampled: 12/20/2013 1257

Date Received: 12/23/2013 0900

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

Lab File ID:

S0101010.D

Dilution:

Prep Batch:

180-93531

Initial Weight/Volume:

20

01/01/2014 1128

Final Weight/Volume:

30.0 g 2.0 mL

Analysis Date: Prep Date:

12/31/2013 0340

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		140	J	30	300
Benzo[a]anthracene		470		38	300
Benzo[b]fluoranthene		800		47	300
Benzo[k]fluoranthene		250	J	61	300
Benzo[g,h,i]perylene		690		30	300
Benzo[a]pyrene		620		30	300
Chrysene		710		36	300
Dibenz(a,h)anthracene		150	J	34	300
Fluoranthene		1000		32	300
Fluorene		43	J	40	300
ndeno[1,2,3-cd]pyrene		540		31	300
Phenanthrene		260	J	48	300
Pyrene		710		31	300
Acenaphthene		ND		29	300
Acenaphthylene		120	J	35	300
Naphthalene		ND		26	300
2-Methylnaphthalene		ND		27	300
Surrogate		%Rec	Qualifier	Accepta	nce Limits
Nitrobenzene-d5 (Surr)	***************************************	84	Ď	27 - 110	
2-Fluorobiphenyl		98	D	28 - 108	
Terphenyl-d14 (Surr)		84	D	21 - 130	

Job Number: 180-28384-1

Client Sample ID:

SD-189-0-1

Lab Sample ID:

180-28384-27

Client Matrix:

Sediment

% Moisture:

9.4

Date Sampled: 12/20/2013 1250

Date Received: 12/23/2013 0900

#### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

21 - 130

71

Prep Method:

3541

Prep Batch:

Lab File ID:

S0101011.D

Dilution:

20

180-93531

Initial Weight/Volume: Final Weight/Volume:

30.0 g

Analysis Date:

01/01/2014 1155 12/31/2013 0340

2.0 mL 2 uL

Prep Date:	12/31/2013 0340		Injec	tion Volume:	2 uL
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene	***************************************	590		29	300
Benzo[a]anthracene		3000		37	300
Benzo[b]fluoranthen	e	3900		46	300
Benzo[k]fluoranthen	e	1500		60	300
Benzo[g,h,i]perylene	•	2800		29	300
Benzo[a]pyrene		3300		29	300
Chrysene		3700		35	300
Dibenz(a,h)anthrace	ne	680		33	300
Fluoranthene		8100		32	300
Fluorene		180	J	39	300
Indeno[1,2,3-cd]pyre	ene	2600		30	300
Phenanthrene		2700		47	300
Pyrene		4500		30	300
Acenaphthene		130	J	28	300
Acenaphthylene		160	J	34	300
Naphthalene		66	J	25	300
2-Methylnaphthalene	e	ND		26	300
Surrogate		%Rec	Qualifier	Accep	otance Limits
Nitrobenzene-d5 (Su	ırr)	74	D	27 - 1	10
2-Fluorobiphenyl		94	D	28 - 1	08

D

79

Terphenyl-d14 (Surr)

Client Sample ID:

SD-190-0-1

Lab Sample ID:

180-28384-28

Client Matrix:

Sediment

% Moisture:

30.2

Date Sampled: 12/20/2013 1445

Date Received: 12/23/2013 0900

8270D LL Semivolatile	Organic (	Compounds	by GC/	MS - Low Level
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Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

Prep Batch:

180-93531

Lab File ID:

S0101012.D

Dilution:

20

Initial Weight/Volume: Final Weight/Volume: 30.0 g

Analysis Date:

01/01/2014 1222

2.0 mL

Prep Date:

12/31/2013 0340

Injection Volume:

12/31/2013 0340			injection voicine.			
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
Anthracene		18000	······································	37	380	conserverententententen
Benzo[a]anthracene		49000		48	380	
Benzo[b]fluoranthene		47000		60	380	
Benzo[k]fluoranthene		19000		77	380	
Benzo[g,h,i]perylene		33000		38	380	
Benzo[a]pyrene		43000		38	380	
Chrysene		55000		45	380	
Dibenz(a,h)anthracene		8200		42	380	
Fluoranthene		130000	E	41	380	
Fluorene		6600		50	380	
Indeno[1,2,3-cd]pyrene		30000		39	380	
Phenanthrene		61000		61	380	
Pyrene		74000		39	380	
Acenaphthene		5600		37	380	
Acenaphthylene		870		44	380	
Naphthalene		120	J	33	380	
2-Methylnaphthalene		530		34	380	
Surrogate		%Rec	Qualifier	Accep	tance Limits	
Nitrobenzene-d5 (Surr)		70	D	27 - 1	10	SCOORSCORPAGEORY-MANUEL
2-Fluorobiphenyl		84	D	28 - 10	08	
Terphenyl-d14 (Surr)		89	D	21 - 13	30	

Client Sample ID:

SD-190-0-1

Lab Sample ID:

180-28384-28

Client Matrix:

Sediment

% Moisture:

30.2

Date Sampled: 12/20/2013 1445

Date Received: 12/23/2013 0900

8270D LL Semivolatile Organic Compounds by 6	GC/MS - Low Level
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Analysis Method:

8270D LL

Analysis Batch:

180-93785

Instrument ID:

71

Prep Method:

3541

Prep Batch:

180-93531

Lab File ID:

S0104012.D

Dilution:

Initial Weight/Volume:

30.0 g

Analysis Date:

50 01/04/2014 1338

Run Type:

DL

Final Weight/Volume:

27 - 110

28 - 108

21 - 130

2.0 mL

Prep Date:

12/31/2013 0340

Injection Volume:

2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		18000	•••••••••••••••••••••••••••••••••••••••	93	960
Benzo[a]anthracene		51000		120	960
Benzo[b]fluoranthene		57000		150	960
Benzo[k]fluoranthene		22000		190	960
Benzo[g,h,i]perylene		40000		95	960
Benzo[a]pyrene		48000		96	960
Chrysene		55000		110	960
Dibenz(a,h)anthracene		9100		110	960
Fluoranthene		150000		100	960
Fluorene		6200		130	960
Indeno[1,2,3-cd]pyrene		33000		98	960
Phenanthrene		71000		150	960
Pyrene		84000		97	960
Acenaphthene		5400		92	960
Acenaphthylene		760	J	110	960
Naphthalene		200	J	82	960
2-Methylnaphthalene		540	J	86	960
Surrogate		%Rec	Qualifier	Acceptance Limits	

D

D

D

62

83

87

Nitrobenzene-d5 (Surr)

Terphenyl-d14 (Surr)

2-Fluorobiphenyl

Job Number: 180-28384-1

Client Sample ID:

SD-191-0-1

Lab Sample ID:

180-28384-29

Client Matrix:

Sediment

% Moisture:

30.1

Date Sampled: 12/20/2013 1450

Date Received: 12/23/2013 0900

### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

Prep Batch:

Lab File ID:

S0101013.D

Dilution:

60

73

66

180-93531

Initial Weight/Volume:

30.1 g

20

D

D

D

Final Weight/Volume:

27 - 110

28 - 108

21 - 130

0.5 mL

Analysis Date: Prep Date:

01/01/2014 1248 12/31/2013 0340

Injection Volume:

2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		4200		9.3	95
Benzo[a]anthracene		13000		12	95
Benzo[b]fluoranthene		16000		15	95
Benzo[k]fluoranthene		5600		19	95
Benzo[g,h,i]perylene		9200		9.5	95
Benzo[a]pyrene		13000		9.5	95
Chrysene		17000		11	95
Dibenz(a,h)anthracene		2300		11	95
Fluoranthene		40000	E	10	95
Fluorene		2400		13	95
Indeno[1,2,3-cd]pyrene		8500		9.8	95
Phenanthrene		22000	E	15	95
Pyrene		23000	E	9.6	95
Acenaphthene		2000		9.1	95
Acenaphthylene		440		11	95
Naphthalene		87	J	8.2	95
2-Methylnaphthalene		230		8.5	95
Surrogate		%Rec	Qualifier	Accepta	nce Limits

Nitrobenzene-d5 (Surr)

Terphenyl-d14 (Surr)

2-Fluorobiphenyl

Job Number: 180-28384-1

Client Sample ID:

SD-191-0-1

Lab Sample ID:

180-28384-29

Client Matrix:

Sediment

% Moisture:

30.1

Date Sampled: 12/20/2013 1450

Date Received: 12/23/2013 0900

# 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93945

Instrument ID:

722

Prep Method:

3541

180-93531

Lab File ID:

F0108001.D

Dilution:

100

Prep Batch:

Initial Weight/Volume:

30.1 g

Analysis Date:

01/08/2014 0223

Run Type:

DL

Final Weight/Volume:

0.5 mL

Prep Date:

12/31/2013 0340

Injection Volume:

2 uL

21 - 130

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		5300	•	47	480
Benzo[a]anthracene		20000		60	480
Benzo[b]fluoranthene		27000		75	480
Benzo[k]fluoranthene		11000		96	480
Benzo[g,h,i]perylene		16000		47	480
Benzo[a]pyrene		21000		48	480
Chrysene		25000		57	480
Dibenz(a,h)anthracene		3200		53	480
Fluoranthene		67000		51	480
Fluorene		3100		63	480
Indeno[1,2,3-cd]pyrene		14000		49	480
Phenanthrene		42000		76	480
Pyrene		49000		48	480
Acenaphthene		2800		46	480
Acenaphthylene		230	J	54	480
Naphthalene		130	J	41	480
2-Methylnaphthalene		290	J	43	480
Surrogate		%Rec	Qualifier	Accepta	nce Limits
Nitrobenzene-d5 (Surr)		0	ΧD	27 - 110	
2-Fluorobiphenyl		0	ΧD	28 - 108	3

ΧD

0

Terphenyl-d14 (Surr)

Client: Tetra Tech, Inc. Job Number: 180-28384-1

Client Sample ID:

SD-192-0-1

Lab Sample ID:

180-28384-30

Client Matrix:

Sediment % Moisture:

23.8

Date Sampled: 12/20/2013 1600

Date Received: 12/23/2013 0900

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

Prep Batch:

Lab File ID:

S0101014.D

Dilution:

180-93531

Initial Weight/Volume:

30.2 g

Analysis Date:

20

Final Weight/Volume:

0.5 mL

Prep Date:

01/01/2014 1315 12/31/2013 0340

Injection Volume:

12.01.2010				<b>,</b>			
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL		
Anthracene		89	**************************************	8.5	87		
Benzo[a]anthracene		130		11	87		
Benzo[b]fluoranthene		170		14	87		
Benzo[k]fluoranthene		64	J	18	87		
Benzo[g,h,i]perylene		120		8.7	87		
Benzo[a]pyrene		150		8.7	87		
Chrysene		190		10	87		
Dibenz(a,h)anthracene		29	J	9.7	87		
Fluoranthene		360		9.3	87		
Fluorene		34	J	11	87		
Indeno[1,2,3-cd]pyrene		99		9.0	87		
Phenanthrene		170		14	87		
Pyrene		240		8.8	87		
Acenaphthene		20	J	8.4	87		
Acenaphthylene		89		10	87		
Naphthalene		ND		7.5	87		
2-Methylnaphthalene		ND		7.8	87		
Surrogate		%Rec	Qualifier	Accepta	nce Limits		
Nitrobenzene-d5 (Surr)		52	D	27 - 110		Indonesianosanerere/errecenonmen	
2-Fluorobiphenyl		67	D	28 - 108	ŀ		
Terphenyl-d14 (Surr)		56	D	21 - 130	1		

Job Number: 180-28384-1

Client Sample ID:

SD-193-0-1

Lab Sample ID:

180-28384-31

Client Matrix:

Sediment

% Moisture:

29.2

Date Sampled: 12/20/2013 1610

Date Received: 12/23/2013 0900

### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

180-93531

Lab File ID:

S0101015.D

Dilution:

Prep Batch:

Initial Weight/Volume:

30.1 g

20

Final Weight/Volume:

0.5 mL

Analysis Date: Prep Date:

01/01/2014 1342 12/31/2013 0340

Injection Volume:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		54	J	9.2	94
Benzo[a]anthracene		160		12	94
Benzo[b]fluoranthene		200		15	94
Benzo[k]fluoranthene		87	J	19	94
Benzo[g,h,i]perylene		130		9.3	94
Benzo[a]pyrene		170		9.4	94
Chrysene		200		11	94
Dibenz(a,h)anthracene		33	J	10	94
luoranthene		440		10	94
Fluorene		23	J	12	94
ndeno[1,2,3-cd]pyrene		110		9.7	94
Phenanthrene		160		15	94
Pyrene		230		9.5	94
Acenaphthene		13	J	9.0	94
Acenaphthylene		31	J	11	94
Naphthalene		ND		8.1	94
2-Methylnaphthalene		ND		8.4	94
Surrogate		%Rec	Qualifier	Acceptance Limits	
Nitrobenzene-d5 (Surr)		58	D	27 - 110	
2-Fluorobiphenyl		72	D	28 - 108	
Terphenyl-d14 (Surr)		89	D	21 - 130	

Job Number: 180-28384-1

Client Sample ID:

SD-194-0-1

Lab Sample ID:

180-28384-32

Client Matrix:

Sediment

% Moisture:

15.7

Date Sampled: 12/20/2013 1615

Date Received: 12/23/2013 0900

#### 8270D LL Semivolatile Organic Compounds by GC/MS - Low Level

Analysis Method:

8270D LL

Analysis Batch:

180-93752

Instrument ID:

71

Prep Method:

3541

Lab File ID:

S0101016.D

Dilution:

Prep Batch:

180-93531

Initial Weight/Volume:

30.0 g

Analysis Date:

2-Fluorobiphenyl

Terphenyl-d14 (Surr)

20

01/01/2014 1409

Final Weight/Volume:

28 - 108 21 - 130 0.5 mL

Prep Date:

12/31/2013 0340

Injection Volume:

2 uL

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
Anthracene		43	J	7.7	80
Benzo[a]anthracene		100		9.9	80
Benzo[b]fluoranthene		140		12	80
Benzo[k]fluoranthene		44	J	16	80
Benzo[g,h,i]perylene		94		7.9	80
Benzo[a]pyrene		130		7.9	80
Chrysene		150		9.4	80
Dibenz(a,h)anthracene		26	J	8.8	80
Fluoranthene		240		8.5	80
Fluorene		15	J	10	80
Indeno[1,2,3-cd]pyrene		79	j	8.2	80
Phenanthrene		81		13	80
Pyrene		160		8.0	80
Acenaphthene		9.0	j	7.6	80
Acenaphthylene		48	J	9.1	80
Naphthalene		ND		6.8	80
2-Methylnaphthalene		ND		7.1	80
Surrogate		%Rec	Qualifier	Acceptance Limits	
Nitrobenzene-d5 (Surr)	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	65	D	27 - 110	

D

D

80

66

Job Number: 180-28384-1

Client Sample ID:

SD-163-0-1

Lab Sample ID:

180-28384-1

Client Matrix:

Sediment

% Moisture:

23.2

Date Sampled: 12/18/2013 1155

Date Received: 12/23/2013 0900

8082A Polychlorinated Biphenyls (PCBs) (GC)	8082A Poly	vchlorinated	<b>Biphenyls</b>	(PCBs)	(GC)
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Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

Initial Weight/Volume:

180-93257

30.1 g

Dilution:

5.0

Final Weight/Volume:

1.0 mL

Analysis Date:

12/27/2013 0427

Injection Volume:

1 uL

Prep Date:

12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND		0.40	2.7	
PCB-1221		ND		0.52	2.7	
PCB-1232		ND		0.46	2.7	
PCB-1242		ND		0.44	2.7	
PCB-1248		ND		0.26	2.7	
PCB-1254		ND		0.38	2.7	
PCB-1260		25		0.38	2.7	
Surrogate		%Rec	Qualifier	Accepta	nce Limits	
DCB Decachlorobiphenyl (Surr	•)	103		20 - 150	)	
Tetrachloro-m-xylene		102		30 - 150	)	

Job Number: 180-28384-1

Client Sample ID:

SD-164-0-1

Lab Sample ID:

180-28384-2

Client Matrix: Sediment % Moisture:

21.2

Date Sampled: 12/18/2013 1210

Date Received: 12/23/2013 0900

8082A Polychlorinated Biphenyls (PCBs) (GC	8082A Pol	ychlorinated	Biphenyls	(PCBs)	(GC)
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Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

Prep Batch:

180-93257

Initial Weight/Volume:

30.0 g

Dilution:

3541

Final Weight/Volume:

Analysis Date:

5.0

Injection Volume:

1.0 mL 1 uL

Prep Date:

12/27/2013 0458 12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		0.39	2.6
PCB-1221		ND		0.50	2.6
PCB-1232		ND		0.45	2.6
PCB-1242		ND		0.43	2.6
PCB-1248		ND		0.25	2.6
PCB-1254		ND		0.38	2.6
PCB-1260		89		0.38	2.6

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	180	X	20 - 150
Tetrachloro-m-xylene	104		30 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-165-0-1

Lab Sample ID:

180-28384-3

Client Matrix:

Sediment

% Moisture:

23.8

Date Sampled: 12/18/2013 1235

Date Received: 12/23/2013 0900

00004	District Landau at a st	Diahaanda	(DCD-) (CC)
BUBZA	Polychlorinated	Bibnenvis	(PUBS) (GU)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

30.1 g

Dilution:

5.0

Final Weight/Volume:

Analysis Date:

Injection Volume:

1.0 mL 1 uL

Prep Date:

12/27/2013 0530 12/26/2013 0340

Result Type:

PRIMARY

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND	······································	0.41	2.7
PCB-1221		ND		0.52	2.7
PCB-1232		ND		0.47	2.7
PCB-1242		ND		0.44	2.7
PCB-1248		ND		0.26	2.7
PCB-1254		ND		0.39	2.7
PCB-1260		99		0.39	2.7
Surrogate		%Rec	Qualifier	Accepta	ance Limits

20 - 150 30 - 150

Job Number: 180-28384-1 Client: Tetra Tech, Inc.

Client Sample ID:

SD-166-0-1

Lab Sample ID:

180-28384-4

Client Matrix:

Sediment

% Moisture:

20.0

Date Sampled: 12/18/2013 1305

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

30.0 g

Dilution:

5.0

Final Weight/Volume:

Analysis Date:

Injection Volume:

1.0 mL 1 uL

Prep Date:

12/27/2013 0704 12/26/2013 0340

Result Type:

PRIMARY

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND		0.39	2.6	
PCB-1221		ND		0.50	2.6	
PCB-1232		ND		0.45	2.6	
PCB-1242		ND		0.42	2.6	
PCB-1248		ND		0.25	2.6	
PCB-1254		ND		0.37	2.6	
PCB-1260		34		0.37	2.6	
Surrogate		%Rec	Qualifier	Accepta	ince Limits	

73 88 20 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-167-0-1

Lab Sample ID:

180-28384-5

Client Matrix:

Sediment

% Moisture:

22.4

Date Sampled: 12/18/2013 1510

Date Received: 12/23/2013 0900

8082A Polychlorinated Biphenyls (PCBs) (GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

Dilution:

30.0 g

5.0

Final Weight/Volume:

1.0 mL

Analysis Date:

Injection Volume:

1 uL

Prep Date:

12/27/2013 0736 12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016	•••••••••••••••••••••••••••••••••••••••	ND		0.40	2.7	•
PCB-1221		ND		0.51	2.7	
PCB-1232		ND		0.46	2.7	
PCB-1242		ND		0.44	2.7	
PCB-1248		ND		0.25	2.7	
PCB-1254		ND		0.38	2.7	
PCB-1260		19		0.38	2.7	
Surrogate		%Rec	Qualifier	Accepta	ance Limits	
DCB Decachlorobiphenyl (Surr	•)	69		20 - 15	0	
Tetrachloro-m-xylene		69		30 - 15	0	

Job Number: 180-28384-1

Client Sample ID:

SD-168-0-1

Lab Sample ID:

180-28384-6

Client Matrix:

Sediment

% Moisture:

21.3

Date Sampled: 12/18/2013 1545

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

30.0 g

Dilution:

5.0

Final Weight/Volume:

1.0 mL

Analysis Date:

12/27/2013 0807

Injection Volume:

1 uL

Prep Date:

12/26/2013 0340

Result Type:

PCB-1016 ND PCB-1221 ND PCB-1232 ND PCB-1242 ND	0.39 0.50 0.45 0.43	2.6 2.6 2.6	
PCB-1232 ND PCB-1242 ND	0.45	2.6	
PCB-1242 ND	- · · · · · · · · · · · · · · · · · · ·	—· <del>-</del>	
	0.43	2.6	
Den 1919	0.43	2.6	
PCB-1248 ND	0.25	2.6	
PCB-1254 100	0.38	2.6	
PCB-1260 84	0.38	2.6	

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	131	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	20 - 150
Tetrachloro-m-xvlene	102		30 - 150

Client: Tetra Tech, Inc.

Job Number: 180-28384-1

Client Sample ID:

SD-169-0-1

Lab Sample ID:

180-28384-7

Client Matrix:

Sediment

% Moisture:

12.4

Date Sampled: 12/18/2013 1610

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

30.1 g

Dilution:

Final Weight/Volume:

Analysis Date:

5.0

1.0 mL

Prep Date:

12/27/2013 0839 12/26/2013 0340

Injection Volume: Result Type:

1 uL PRIMARY

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		0.35	2.4
PCB-1221		ND		0.45	2.4
PCB-1232		ND		0.41	2.4
PCB-1242		ND		0.39	2.4
PCB-1248		ND		0.22	2.4
PCB-1254		29		0.34	2.4
PCB-1260		29		0.34	2.4

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	74		20 - 150
Tetrachloro-m-xvlene	76		30 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-170-0-1

Lab Sample ID:

180-28384-8

Client Matrix:

Sediment

% Moisture:

21.5

Date Sampled: 12/20/2013 1515

Date Received: 12/23/2013 0900

	8082A Polychlorinated	Biphenyls (	(PCBs) (GC	)
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Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

Prep Batch:

180-93257

Initial Weight/Volume:

3541

Final Weight/Volume:

30.1 g

Dilution:

1.0 mL

Analysis Date:

5.0

Injection Volume:

1 uL

Prep Date:

12/27/2013 0910 12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		0.39	2.6
PCB-1221		ND		0.51	2.6
PCB-1232		ND		0.45	2.6
PCB-1242		ND		0.43	2.6
PCB-1248		ND		0.25	2.6
PCB-1254		ND		0.38	2.6
PCB-1260		2.1	J	0.38	2.6
Surrogate		%Rec	Qualifier	Accepta	nce Limits
DCB Decachlorobiphenyl (Surr	r)	64	***************************************	20 - 150	)

Job Number: 180-28384-1

Client Sample ID:

SD-171-0-1

Lab Sample ID:

180-28384-9

Client Matrix:

Sediment

% Moisture:

66.8

Date Sampled: 12/19/2013 1045

Date Received: 12/23/2013 0900

8082A Polychlorinated Biphenyls (PCBs) (GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

Initial Weight/Volume:

180-93257

30.0 g

Dilution:

1000

Final Weight/Volume:

1.0 mL

Analysis Date:

Injection Volume:

1 uL

Prep Date:

12/27/2013 1909 12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016	······································	ND	······································	190	1300	
PCB-1221		ND		240	1300	
PCB-1232		ND		210	1300	
PCB-1242		ND		200	1300	
PCB-1248		ND		120	1300	
PCB-1254		ND		180	1300	
PCB-1260		150000		180	1300	
Surrogate		%Rec	Qualifier	Accepta	ance Limits	
DCB Decachlorobiphenyl (Suri	r)	0	ΧD	20 - 150	)	***************************************
Tetrachloro-m-xylene		0	ΧD	30 - 150	)	

Job Number: 180-28384-1

Client Sample ID:

SD-172-0-3

Lab Sample ID:

180-28384-10

Client Matrix:

Sediment

% Moisture:

46.9

Date Sampled: 12/19/2013 1110

Date Received: 12/23/2013 0900

8082A Polychlorinated Biphenyls (PCBs) (GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

0

180-93257

Initial Weight/Volume:

30 - 150

Dilution:

30.0 g

1000

Final Weight/Volume:

1.0 mL

Analysis Date:

Injection Volume:

1 uL

Prep Date:

Tetrachloro-m-xylene

12/27/2013 1940 12/26/2013 0340

Result Type:

**PRIMARY** 

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND	***************************************	120	780
PCB-1221		ND		150	780
PCB-1232		ND		130	780
PCB-1242		ND		130	780
PCB-1248		ND		74	780
PCB-1254		ND		110	780
PCB-1260		54000		110	780
Surrogate		%Rec	Qualifier	Accepta	ince Limits
DCB Decachlorobiphenyl (Surr)	········	0	ΧD	20 - 150	)

ΧD

Job Number: 180-28384-1

Client Sample ID:

SD-173-0-3

Lab Sample ID:

180-28384-11

Client Matrix:

Sediment

% Moisture:

38.0

Date Sampled: 12/19/2013 1130

Date Received: 12/23/2013 0900

8082A Pol	vchlorinated	Biphenyls	(PCBs)	(GC)
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Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

30.0 g

Dilution:

5000

Final Weight/Volume:

1.0 mL

Analysis Date:

Injection Volume:

1 uL

Prep Date:

12/27/2013 2012 12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND		500	3400	
PCB-1221		ND		640	3400	
PCB-1232		ND		580	3400	
PCB-1242		ND		550	3400	
PCB-1248		ND		320	3400	
PCB-1254		ND		480	3400	
PCB-1260		220000		480	3400	
Surrogate		%Rec	Qualifier	Accept	ance Limits	
DCB Decachlorobiphenyl (Sur	·r)	0	ΧD	20 - 15	0	
Tetrachloro-m-xylene		0	ΧD	30 - 15	0	

Job Number: 180-28384-1

Client Sample ID:

SD-174-0-1

Lab Sample ID:

180-28384-12

Client Matrix:

Sediment

% Moisture:

16.6

Date Sampled: 12/19/2013 1340

Date Received: 12/23/2013 0900

8082A Pc	lychlorinated	Biphenvis	(PCBs)	(GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

Initial Weight/Volume:

180-93257

30.1 g

Dilution:

5.0

Final Weight/Volume:

1.0 mL

Analysis Date:

Injection Volume:

1 uL

Prep Date:

12/27/2013 1426 12/26/2013 0340

Result Type:

A -1.4	D 1440	D = 14 ( = 44=)	Overlife	MDI	RL	
Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL		••••
PCB-1016		ND		0.37	2.5	
PCB-1221		ND		0.48	2.5	
PCB-1232		ND		0.43	2.5	
PCB-1242		ND		0.41	2.5	
PCB-1248		ND		0.24	2.5	
PCB-1254		150		0.35	2.5	
PCB-1260		350		0.35	2.5	
Surrogate		%Rec	Qualifier	Accepta	nce Limits	
DCB Decachlorobiphenyl (Sur	r)	131		20 - 150	)	***********************
Tetrachloro-m-xylene	•	96		30 - 150	)	

Job Number: 180-28384-1

Client Sample ID:

SD-175-0-1

Lab Sample ID:

180-28384-13

Client Matrix:

Sediment

% Moisture:

30.0

Date Sampled: 12/19/2013 1430

Date Received: 12/23/2013 0900

8082A Polychlorinated Biphenyls (PCBs) (GC)
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Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

Final Weight/Volume:

30.0 g

Dilution:

50

1.0 mL

Analysis Date:

Injection Volume:

1 uL

Prep Date:

12/28/2013 0816 12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		4.4	30
PCB-1221		ND		5.7	30
PCB-1232		ND		5.1	30
PCB-1242		ND		4.8	30
PCB-1248		ND		2.8	30
PCB-1254		410		4.2	30
PCB-1260		1300		4.2	30
Surrogate		%Rec	Qualifier	Acc	ceptance Limits
DCB Decachlorobiphenyl (Sur	r)	180	DX	20 -	- 150
Tetrachloro-m-xylene	•	120	D	30 -	- 150

Job Number: 180-28384-1

Client Sample ID:

SD-176-0-3

Lab Sample ID:

180-28384-14

Client Matrix:

Sediment

% Moisture:

43.3

Date Sampled: 12/19/2013 1505

Date Received: 12/23/2013 0900

8082A F	<b>Polychlorinated</b>	Biphenyls i	(PCBs)	(GC)
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Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

30.0 g

Dilution:

5.0

Final Weight/Volume:

1.0 mL

Analysis Date:

Injection Volume:

1 uL

Prep Date:

12/27/2013 1528 12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		0.55	3.7
PCB-1221		ND		0.70	3.7
PCB-1232		ND		0.63	3.7
PCB-1242		ND		0.60	3.7
PCB-1248		ND		0.35	3.7
PCB-1254		36		0.52	3.7
PCB-1260		88		0.52	3.7
Surrogate		%Rec	Qualifier	Accepta	ance Limits

Surrogate	%Rec	Qualifier	Acceptance
DCB Decachlorobiphenyl (Surr)	88		20 - 150
Tetrachloro-m-xylene	89		30 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-177-0-2

Lab Sample ID:

180-28384-15

Client Matrix:

Sediment

% Moisture:

71.8

Date Sampled: 12/19/2013 1540

Date Received: 12/23/2013 0900

8082A	Polychlorinated	Biphenyls	(PCBs)	(GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

30.2 g

Dilution:

5.0

Final Weight/Volume:

1.0 mL

Analysis Date:

12/27/2013 1600

Injection Volume:

1 uL

Prep Date:

12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND		1.1	7.3	***************************************
PCB-1221		ND		1.4	7.3	
PCB-1232		ND		1.3	7.3	
PCB-1242		ND		1.2	7.3	
PCB-1248		ND		0.69	7.3	
PCB-1254		28		1.0	7.3	
PCB-1260		73		1.0	7.3	
Surrogate		%Rec	Qualifier	Accepta	ince Limits	
DCB Decachlorobiphenyl (Surr	·····	84		20 - 150	)	***************************************
Tetrachloro-m-xylene		93		30 - 150	)	

Job Number: 180-28384-1

Client Sample ID:

SD-178-0-1

Lab Sample ID:

180-28384-16

Client Matrix:

Sediment

% Moisture:

21.4

Date Sampled: 12/19/2013 1600

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

30.1 g

Dilution:

5.0

Final Weight/Volume:

1.0 mL

Analysis Date:

12/27/2013 1631

Injection Volume:

1 uL

Prep Date:

12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		0.39	2.6
PCB-1221		ND		0.50	2.6
PCB-1232		ND		0.45	2.6
PCB-1242		ND		0.43	2.6
PCB-1248		ND		0.25	2.6
PCB-1254		5.3		0.38	2.6
PCB-1260		12		0.38	2.6

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	62	//////////////////////////////////////	20 - 150
Tetrachloro-m-xylene	75		30 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-179-0-1

Lab Sample ID:

180-28384-17

Client Matrix:

Sediment

% Moisture:

30.3

Date Sampled: 12/19/2013 1645

Date Received: 12/23/2013 0900

8082A Polychlorinated	Biphenyls	(PCBs)	(GC)	
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Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

Initial Weight/Volume:

180-93257

30.0 g

Dilution:

5.0

Final Weight/Volume:

1.0 mL

Analysis Date:

Injection Volume:

1 uL

Prep Date:

12/27/2013 1703 12/26/2013 0340

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016	······································	ND		0.44	3.0	***************************************
PCB-1221		ND		0.57	3.0	
PCB-1232		ND		0.51	3.0	
PCB-1242		ND		0.49	3.0	
PCB-1248		ND		0.28	3.0	
PCB-1254		7.6		0.42	3.0	
PCB-1260		22		0.42	3.0	
Surrogate		%Rec	Qualifier	Accepta	ance Limits	
DCB Decachlorobiphenyl (Surr	······································	81		20 - 150	)	
Tetrachloro-m-xylene		94		30 - 150	)	

Job Number: 180-28384-1

Client Sample ID:

SD-180-0-1

Lab Sample ID:

180-28384-18

Client Matrix:

Sediment

% Moisture:

21.3

Date Sampled: 12/19/2013 1625

Date Received: 12/23/2013 0900

8082A Polychlorinated	Biphenvis	(PCBs)	(GC)
OUOLA : Olyomormated	D.p,	·,	ι,

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

Prep Batch:

180-93257

Initial Weight/Volume:

3541

30.0 g

Dilution:

10

Final Weight/Volume:

1.0 mL

Analysis Date:

12/28/2013 0848

Injection Volume:

1 uL

Prep Date:

12/26/2013 0340

Result Type:

**PRIMARY** 

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND		0.79	5.3	***************************************
PCB-1221		ND		1.0	5.3	
PCB-1232		ND		0.91	5.3	
PCB-1242		ND		0.86	5.3	
PCB-1248		ND		0.50	5.3	
PCB-1254		21		0.75	5.3	
PCB-1260		95		0.75	5.3	
Surrogate		%Rec	Qualifier	Accepta	nce Limits	

20 - 150 30 - 150

Client: Tetra Tech, Inc.

Job Number: 180-28384-1

Client Sample ID:

SD-181-0-1

Lab Sample ID:

180-28384-19

Client Matrix:

Sediment

% Moisture:

42.5

Date Sampled: 12/20/2013 1000

Date Received: 12/23/2013 0900

### 8082A Polychlorinated Biphenyls (PCBs) (GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

30.1 g

Dilution:

Final Weight/Volume:

Analysis Date:

20000

1.0 mL 1 uL

Prep Date:

12/28/2013 1158 12/26/2013 0340

Injection Volume: Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND		2100	14000	
PCB-1221		ND		2800	14000	
PCB-1232		ND		2500	14000	
PCB-1242		ND		2400	14000	
PCB-1248		ND		1400	14000	
PCB-1254		ND		2100	14000	
PCB-1260		420000		2100	14000	
Surrogate		%Rec	Qualifier	Accepta	nce Limits	
DCB Decachlorobiphenyl (Surr	)	0	ΧD	20 - 150		200000000000000000000000000000000000000
Tetrachloro-m-xylene	•	0	ΧD	30 - 150	)	

Job Number: 180-28384-1

Client Sample ID:

SD-182-0-3

Lab Sample ID:

180-28384-20

Client Matrix:

Sediment

% Moisture:

52.1

Date Sampled: 12/20/2013 1020

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93257

Initial Weight/Volume:

Dilution:

20000

30.0 g

Final Weight/Volume: Injection Volume:

1.0 mL

Analysis Date: Prep Date:

12/28/2013 1229 12/26/2013 0340

Result Type:

1 uL PRIMARY

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		2600	17000
PCB-1221		ND		3300	17000
PCB-1232		ND		3000	17000
PCB-1242		ND		2800	17000
PCB-1248		ND		1600	17000
PCB-1254		ND		2500	17000
PCB-1260		780000		2500	17000
Surrogate		%Rec	Qualifier	Accepta	nce Limits
DCB Decachlorobiphenyl (Surr	)	0	ΧD	20 - 150	
Tetrachloro-m-xylene	•	0	ΧD	30 - 150	)

Job Number: 180-28384-1

Client Sample ID:

SD-183-0-1

Lab Sample ID:

180-28384-21

Client Matrix:

Sediment

% Moisture:

69.9

Date Sampled: 12/20/2013 1040

Date Received: 12/23/2013 0900

8082A Pol	ychlorinated	<b>Bipheny</b>	Is (PCBs	) (GC)
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Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Initial Weight/Volume:

Prep Batch:

180-93333

30.0 g

Dilution:

2000

Final Weight/Volume:

1.0 mL

Analysis Date:

12/28/2013 1023

Injection Volume:

1 uL

Prep Date:

12/27/2013 0240

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND ND		410	2800	
PCB-1221		ND		530	2800	
PCB-1232		ND		470	2800	
PCB-1242		ND		450	2800	
PCB-1248		ND		260	2800	
PCB-1254		ND		390	2800	
PCB-1260		40000		390	2800	
Surrogate		%Rec	Qualifier	Accepta	ance Limits	
DCB Decachlorobiphenyl (Surr	)	0	ХD	20 - 150	0	**********
Tetrachloro-m-xylene	,	0	ΧD	30 - 150	0	

Job Number: 180-28384-1 Client: Tetra Tech, Inc.

Client Sample ID:

SD-184-0-2

Lab Sample ID:

180-28384-22

Client Matrix:

Sediment

% Moisture:

32.1

Date Sampled: 12/20/2013 1130

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Initial Weight/Volume:

Prep Batch:

180-93333

30.2 g

Dilution:

10

Final Weight/Volume:

1.0 mL

Analysis Date: Prep Date:

12/27/2013 2321

12/27/2013 0240

Injection Volume: Result Type:

1 uL **PRIMARY** 

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		0.91	6.1
PCB-1221		ND		1.2	6.1
PCB-1232		ND		1.0	6.1
PCB-1242		ND		0.99	6.1
PCB-1248		ND		0.58	6.1
PCB-1254		410		0.87	6.1
PCB-1260		1300		0.87	6.1

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	105	***************************************	20 - 150
Tetrachloro-m-xviene	89		30 - 150

Job Number: 180-28384-1 Client: Tetra Tech, Inc.

Client Sample ID:

SD-185-0-1

Lab Sample ID:

180-28384-23

Client Matrix:

Sediment

% Moisture:

25.4

Date Sampled: 12/20/2013 1200

Date Received: 12/23/2013 0900

8082A Polychlorinated Biphenyls (PCBs) (GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93333

Initial Weight/Volume:

30.0 g

Dilution:

10

Final Weight/Volume:

Analysis Date:

Injection Volume:

1.0 mL 1 uL

Prep Date:

12/27/2013 2352 12/27/2013 0240

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		0.83	5.6
PCB-1221		ND		1.1	5.6
PCB-1232		ND		0.96	5.6
PCB-1242		ND		0.91	5.6
PCB-1248		ND		0.53	5.6
PCB-1254		190		0.79	5.6
PCB-1260		550		0.79	5.6

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	103		20 - 150
Tetrachloro-m-xylene	123		30 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-186-0-1

Lab Sample ID:

180-28384-24

Client Matrix:

Sediment

% Moisture:

25.8

Date Sampled: 12/20/2013 1220

Date Received: 12/23/2013 0900

8082A Pol	ychlorinated	Biphenvls	(PCBs)	(GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93333

Initial Weight/Volume:

30.0 g

Dilution:

Final Weight/Volume:

10

1.0 mL

Analysis Date:

12/28/2013 0127

Injection Volume:

1 uL

Prep Date:

12/27/2013 0240

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND	······································	0.84	5.6
PCB-1221		ND		1.1	5.6
PCB-1232		ND		0.96	5.6
PCB-1242		ND		0.91	5.6
PCB-1248		ND		0.53	5.6
PCB-1254		38		0.80	5.6
PCB-1260		60		0.80	5.6

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	64	//////////////////////////////////////	20 - 150
Tetrachloro-m-xvlene	87		30 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-187-0-1

Lab Sample ID:

180-28384-25

Client Matrix:

Sediment

% Moisture:

20.3

Date Sampled: 12/20/2013 1245

Date Received: 12/23/2013 0900

2022	Polychlorinated	Rinhenvis	(PCRs)	(GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

Initial Weight/Volume:

Dilution:

180-93333

30.1 g

10

Final Weight/Volume:

1.0 mL

Analysis Date:

12/28/2013 0158

Injection Volume:

1 uL

Prep Date:

12/27/2013 0240

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016	······································	ND		0.78	5.2
PCB-1221		ND		0.99	5.2
PCB-1232		ND		0.89	5.2
PCB-1242		ND		0.85	5.2
PCB-1248		ND		0.49	5.2
PCB-1254	•	29		0.74	5.2
PCB-1260		92		0.74	5.2
Surrogate		%Rec	Qualifier	Accepta	ance Limits
DCB Decachlorobinhenyl (Surr	\ \	65		20 - 150	)

Job Number: 180-28384-1

Client Sample ID:

SD-188-0-1

Lab Sample ID:

180-28384-26

Client Matrix:

Sediment

% Moisture:

11.7

Date Sampled: 12/20/2013 1257

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Initial Weight/Volume:

Dilution:

Prep Batch:

180-93333

Final Weight/Volume:

30.0 g

Analysis Date:

10

1.0 mL

Prep Date:

12/28/2013 0230 12/27/2013 0240

Injection Volume: Result Type:

1 uL PRIMARY

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		0.70	4.7
PCB-1221		ND		0.90	4.7
PCB-1232		ND		0.81	4.7
PCB-1242		ND		0.77	4.7
PCB-1248		ND		0.45	4.7
PCB-1254		15		0.67	4.7
PCB-1260		49		0.67	4.7

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	64		20 - 150
Tetrachloro-m-xvlene	77		30 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-189-0-1

Lab Sample ID:

180-28384-27

Client Matrix:

Sediment

% Moisture:

9.4

Date Sampled: 12/20/2013 1250

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93333

Initial Weight/Volume:

30.1 g

Dilution:

Final Weight/Volume:

Analysis Date:

10

1.0 mL

12/28/2013 0301

Injection Volume:

1 uL

Prep Date:

12/27/2013 0240

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND		0.68	4.6
PCB-1221		ND		0.87	4.6
PCB-1232		ND		0.78	4.6
PCB-1242		ND		0.75	4.6
PCB-1248		ND		0.43	4.6
PCB-1254		11		0.65	4.6
PCB-1260		82		0.65	4.6

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	70		20 - 150
Tetrachloro-m-xylene	82		30 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-190-0-1

Lab Sample ID:

180-28384-28

Client Matrix:

Sediment

% Moisture:

30.2

Date Sampled: 12/20/2013 1445

Date Received: 12/23/2013 0900

### 8082A Polychlorinated Biphenyls (PCBs) (GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Initial Weight/Volume:

30.1 g

Prep Batch:

180-93333

Dilution:

10

Final Weight/Volume:

1.0 mL

Analysis Date:

12/28/2013 0333

Injection Volume:

1 uL

Prep Date:

12/27/2013 0240

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016	·	ND		0.88	5.9	**********
PCB-1221		ND		1.1	5.9	
PCB-1232		ND		1.0	5.9	
PCB-1242		ND		0.97	5.9	
PCB-1248		ND		0.56	5.9	
PCB-1254		53		0.85	5.9	
PCB-1260		46		0.85	5.9	
Surrogate		%Rec	Qualifier	Accepta	nce Limits	
DCB Decachlorobiphenyl (Surr)	······································	53	***************************************	20 - 150	)	der der dende de
Tetrachloro-m-xylene		81		30 - 150	)	

Job Number: 180-28384-1 Client: Tetra Tech, Inc.

Client Sample ID:

SD-191-0-1

Lab Sample ID:

180-28384-29

Client Matrix:

Sediment

% Moisture:

30.1

Date Sampled: 12/20/2013 1450

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93333

Initial Weight/Volume:

30.0 g

Dilution:

Final Weight/Volume:

Analysis Date:

10

Injection Volume:

1.0 mL 1 uL

Prep Date:

12/28/2013 0404 12/27/2013 0240

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND		0.89	6.0	***************************************
PCB-1221		ND		1.1	6.0	
PCB-1232		ND		1.0	6.0	
PCB-1242		ND		0.97	6.0	
PCB-1248		ND		0.56	6.0	
PCB-1254		350		0.85	6.0	
PCB-1260		530		0.85	6.0	
Surrogate		%Rec	Qualifier	Accepta	nce Limits	

Surrogate	70Rec	Quaimer	Acceptance
DCB Decachlorobiphenyl (Surr)	89		20 - 150
Tetrachloro-m-xylene	82		30 - 150

Job Number: 180-28384-1 Client: Tetra Tech, Inc.

Client Sample ID:

SD-192-0-1

Lab Sample ID:

180-28384-30

Client Matrix:

Sediment

% Moisture:

23.8

Date Sampled: 12/20/2013 1600

Date Received: 12/23/2013 0900

8082A Pol	ychlorinated	<b>Biphenvis</b>	(PCBs)	(GC)
000271 01	, o o a . c . a	D.p.io.iyio	(. 000,	,,

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

180-93333

Initial Weight/Volume:

Dilution:

Prep Batch:

30.0 g

10

Final Weight/Volume:

1.0 mL

Analysis Date: Prep Date:

12/28/2013 0436

12/27/2013 0240

Injection Volume: 1 uL

Result Type:

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND		0.81	5.5	
PCB-1221		ND		1.0	5.5	
PCB-1232		ND		0.94	5.5	
PCB-1242		ND		0.89	5.5	
PCB-1248		ND		0.52	5.5	
PCB-1254		13		0.78	5.5	
PCB-1260		20		0.78	5.5	

Surrogate	%Rec	Qualifier	Acceptance Limits
DCB Decachlorobiphenyl (Surr)	54		20 - 150
Tetrachloro-m-xylene	65		30 - 150

Job Number: 180-28384-1 Client: Tetra Tech, Inc.

Client Sample ID:

SD-193-0-1

Lab Sample ID:

180-28384-31

Client Matrix:

Sediment

% Moisture:

29.2

Date Sampled: 12/20/2013 1610

Date Received: 12/23/2013 0900

### 8082A Polychlorinated Biphenyls (PCBs) (GC)

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93333

Initial Weight/Volume:

30.0 g

10

Dilution:

Final Weight/Volume:

Analysis Date:

Injection Volume:

1.0 mL 1 uL

Prep Date:

12/28/2013 0507 12/27/2013 0240

Result Type:

**PRIMARY** 

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL
PCB-1016		ND	**************************************	0.88	5.9
PCB-1221		ND		1.1	5.9
PCB-1232		ND		1.0	5.9
PCB-1242		ND		0.96	5.9
PCB-1248		ND		0.56	5.9
PCB-1254		ND		0.84	5.9
PCB-1260		17		0.84	5.9
Surrogate		%Rec	Qualifier	Accepta	ance Limits
DCB December high and /Curr	\	60		20 150	````

20 - 150 30 - 150

Job Number: 180-28384-1

Client Sample ID:

SD-194-0-1

Lab Sample ID:

180-28384-32

Client Matrix:

Sediment

% Moisture:

15.7

Date Sampled: 12/20/2013 1615

Date Received: 12/23/2013 0900

Analysis Method:

8082A

Analysis Batch:

180-93552

Instrument ID:

GC10

Prep Method:

3541

Prep Batch:

180-93333

Initial Weight/Volume:

30.1 g

Dilution:

10

Final Weight/Volume:

1.0 mL

Analysis Date:

Injection Volume:

1 uL

Prep Date:

12/28/2013 0539 12/27/2013 0240

Result Type:

**PRIMARY** 

Analyte	DryWt Corrected: Y	Result (ug/Kg)	Qualifier	MDL	RL	
PCB-1016		ND	······································	0.73	4.9	
PCB-1221		ND		0.94	4.9	
PCB-1232		ND		0.84	4.9	
PCB-1242		ND		0.80	4.9	
PCB-1248		ND		0.47	4.9	
PCB-1254		ND		0.70	4.9	
PCB-1260		16		0.70	4.9	
Surrogate		%Rec	Qualifier	Accepta	nce Limits	

DCB Decachlorobiphenyl (Surr)
Tetrachloro-m-xylene

20 - 150 30 - 150

# Appendix C

Support Documentation

# Chain of Custody Record

**TestAmerica** 

The Leacher in Envisoration is sinke N - None
O - Ashao2
P - Na2O45
Q - Na2O45
Q - Na2S03
R - Na2S2S03
S - H2SO4
T - TSP Dodecatrydrate
U - Acetone
W - InfoA
W - Inf - 5
Z - other (specify) Special Instructions/Note: Detertine: 12/20/13 (430 Felex 12xc05483 Sample Disposal ( A fee may be assessed if samples are retained longer than 1 month)

Return To Client Disposal By Lab Monte For Monte Special Instructions/QC Requirements. 180-15270-4013.1 A - HCL
B - NaOH
C - Zn Acetate
D - Nüric Acid
E - NaHSO4
F - Mod - Marso4
F - Mod - Andahor
H - Ascorbic Acid
I - Ice
J - DI Water
K - EDTA
L - EDA Page 1 of 3 180-28384 Chain of Custody 8007 1903 9380 Aethod of Shipment FEDEY **Analysis Requested** Cooler Temperature(s) C and Other Remarks: Fred Kolber & Tt-Gen Bowen, Debra one: 524, 3029 Gebra. bowen@testamericainc.com (eceived by: PLL रकाव SHAG ×××× XXXX Company Tetra Teck Company Matrix (w-water, g-solid, O-wastaskid, Company ŋ Sample (\* Type (\* C=comp, c=c) G=grab) sr-ris Radiological 301. 524. 3029 1730 1130 1305 9191 12/19/13 1045 Sample 735 1510 1110 1710 25.57 2/20/13/15/15 12/18/13 1188 21 days (AT Requested (days): Unknown 14/20/13
Date/Time Due Date Requested: Sample Date Project#: 18012067 SSOW#: PO#: 1097251 ₩OM Poison B Storn Draw Sediment Invest. Skin Imtant Deliverable Requested: I, II, III, IV, Other (specify) Custody Seal No. fiddle River Complex, Middle River MD Flammable Address: 2171 West Park Court Suite E Possible Hazard Identification Eric.Samuels@tetratech.com 84-173-0-3 SD-164-0-SD-[63-0-50-577-0-3 SD-165-0-1-0-691-05 SD-166-0-1 50-169-0-1 sp-168-0-1 1-0-061-45 SD-171-0-Empty Kit Relinquished by: Custody Seals Intact:

Δ Yes Δ No Client Information Sample Identification Non-Hazard Relinquished by. company: etra Tech, Inc. Stone Mountain inc Samuels elinquished by: elinquished by: State, Zip: GA, 30087

# Chain of Custody Record

**TestAmerica** 

THE LEADER IN CRUMONNESS AL TESTING N - None
O - Ashado
P - Na204S
Q - Na2SSO3
R - Na2SSSO3
R - Na2SSSO3
S - H2SO4
T - TSP Dodecahydrate
U - Acetone
V - MCAA
W - ph 4-5
Z - other (specify) Special Instructions/Note: 1125cosy83 Carrier Tracking No(s): 000 No: 180-15270-4013.2 | 180-15270-4013.2 | Page: 03 Sample Disposal ( A fee may be assessed if samples are retained longer than 1 month)

Return To Client Disposal By Lab Archive For Monti D - Nitric Acid E - NaHSO4 F - MeOH G - Amchlor H - Ascorbic Acid Date/Time: Date/19 1830 ]--loe J--DI Water K-EDTA L-EDA Archive For Method of Shipment **Analysis Requested** Sooler Temperature(s) °C and Other Remarks: Special Instructions/QC Requirements Lab PM: Bowen, Debra E-Mail: debra.bowen@testamericainc.com × × × ×× Company
Tetra Teh
Company Matrix (w-water, 8-soler, O-westeloli, Сотрапу 4 Sample Type (C=comp, G=grab) Radiological Sampler Fred Kothern 19:30 Phone: 301. 578.3014 (340 15.40 1000 Sample Time 1430 1505 1600 1645 001 oha/ 1130 Kyp It Date: AT Requested (days); Unknown Oue Date Requested; 12/19/13 Date/Tipe: 13 Sample Date E1/07/13 Project #. 18012067 SSOW#: PO ≢: 1097251 WO #: Date/Time: Poison B STORM Drain Sedianat Lavest. Skin Irritant Xon-Hazard Flammable Skin Infte Deliverable Requested: I, II, IV, Other (specify) Custody Seal No. Middle River Complex, Middle River MD 2171 West Park Court Suite E ossible Hazard Identification E-0-2, 1943 Eric. Samuels@tetratech.com SD-199-0-SD-184-0-2 50-182-0-3 50-183-0-1 50-174-0-1 1-0-081-0S 50-175-0-1 1-0-181-05 Empty Kit Relinquished by Client Information Custody Seals Intact:
A Yes A No Sample Identification Stone Mountain ompany: etra Tech, Inc. iric Samuels elinguished by: State, Zip: GA, 30087 Relinquished by:

# Chain of Custody Record

TestAmerica

1830 Frais N - Norse
O - AsNaGO
P - NazOGS
Q - NazSOS
R - NazSSSOS
S - 12SO4
T - TSP Dodecabydra
U - Acetone
V - MCAA
W - th 4-5
Z - other (specify) Special Instructions/Note: 112xcost83 Months Sample Disposal ( A fee may be assessed if samples are retained longer than 1 month)

Return To Client Disposal By Lab Archive For Month
Special Instructions/QC Requirements: COC No: 180-15270-4013.3 Pago: D - Nitric Acid E - NaHSO4 F - MeCH G - Amchlor H - Ascorbic Acid Page 3 of 🕏 I-loe J-DI Water K-EDTA L-EDA Date/Time: Jaly 1830 8007 1403 7380 Finex Method of Shipment **Analysis Requested** Cooler Temperature(s) °C and Other Remarks: Lab PM: Bowen, Debra E-Mai: debra. bowen@testamericainc.com PAHE, PCBS, × × X ×× Tela Tech Matrix (w-water, S-roole, O-wastevol, Company G Sample Type (C=comp, G=grab) Radiological hone 301.574.3019 b Fred Korberr 1830 1600 1445 1450 Sample なな 052) 1610 0021 81/02/EI वरत) 342 1615 [AT Requested (days); Unknown Due Dafe Requested: Date/Time: Sample Date Project#: 18012067 PO#: 1097251 \*MOS ₩Ow )ate/∏ime Poison B STOLA DAALH SEULMENT TAVEST. Skin Imitant Possible Hazard Identification

(Mon-Hazard — Fiammable — Skin Imi
Deliverable Requested: I, III, IV, Other (specify) Custody Seal No. <sup>ইতি</sup>ভব Nane: Middle River Complex, Middle River MD 2171 West Park Court Suite E Eric. Samuels@tetratech.com SD-192 - 0-1 50-193-0-1 SD-189-0-20-190-0-50-191-0-Empty Kit Relinquished by: SD-188-0-20-186-0-SD-187-6-1 SD-194-0-1 \$D-185-0-1 Client Information Custody Seals Intact:

Δ Yes Δ No Sample Identification Stone Mountain etra Tech, Inc. Client Contact Eric Samuels delinquished by: Relinquished by: State, Zip: GA, 30087 elinquished by:

### Login Sample Receipt Checklist

Client: Tetra Tech, Inc.

Job Number: 180-28384-1

Login Number: 28384

List Source: TestAmerica Pittsburgh

List Number: 1 Creator: Kovitch, Christina M

Question	Answer Co	mment
Radioactivity wasn't checked or is = background as measured by a survey meter.</td <td>True</td> <td></td>	True	
The cooler's custody seal, if present, is intact.	True	
Sample custody seals, if present, are intact.	True	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
Is the Field Sampler's name present on COC?	True	
There are no discrepancies between the containers received and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	N/A	

### **DATA REPORTING QUALIFIERS**

Client: Tetra Tech, Inc.

Job Number: 180-28384-1

Lab Section	Qualifier	Description
GC/MS Semi VOA		
	*	ISTD response or retention time outside acceptable limits
	F1	MS and/or MSD Recovery exceeds the control limits
	4	MS, MSD: The analyte present in the original sample is greater than 4 times the matrix spike concentration; therefore, control limits are not applicable.
	F2	MS/MSD RPD exceeds control 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.
	X	Surrogate is outside control limits
	D	Surrogate or matrix spike recoveries were not obtained because the extract was diluted for analysis; also compounds analyzed at a dilution may be flagged with a D.
GC Semi VOA		
	4	MS, MSD: The analyte present in the original sample is greater than 4 times the matrix spike concentration; therefore, control limits are not applicable.
	F	MS/MSD Recovery and/or RPD exceeds the control limits
	J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.
	X	Surrogate is outside control limits
	D	Surrogate or matrix spike recoveries were not obtained because the extract was diluted for analysis; also compounds analyzed at a dilution may be flagged with a D.

### **CASE NARRATIVE**

Client: Tetra Tech, Inc.

**Project: Middle River Complex, Middle River MD** 

Report Number: 180-28384-1

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.

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.

### **RECEIPT**

The samples were received on 12/23/2013; the samples arrived in good condition and on ice. The temperature of the cooler at receipt was 1.8° C.

### SEMIVOLATILE ORGANIC COMPOUNDS (GC/MS)

The samples and associated MS/MSD were analyzed at a dilution due to the abundance of target analytes and/or matrix interference. The reporting limits have been adjusted accordingly.

The surrogates of several samples are considered diluted out.

Several analytes failed the recovery criteria for the MS and MSD of sample SD-163-0-1 (180-28384-1).

Several analytes failed the recovery criteria for the MS and/or MSD of sample SD-185-0-1 (180-28384-23). The presence of the '4' qualifier in the data indicates analytes where the concentration in the unspiked sample exceeded four times the spiking amount. Several analytes also exceeded the relative percent difference (RPD) limits.

Internal standard responses were outside of acceptance limits for several samples.

### POLYCHLORINATED BIPHENYLS (PCBs)

The samples and associated MS/MSD were analyzed at a dilution due to the abundance of target analytes and/or matrix interference. The reporting limits have been adjusted accordingly.

DCB Decachlorobiphenyl (Surr) failed the surrogate recovery criteria high for sample SD-164-0-1 (180-28384-2). Tetrachloro-m-xylene recovered within acceptable QC limits. The data are reported.

The surrogates of several samples are considered diluted out.

PCB-1260 failed the recovery criteria for the MS/MSD of sample SD-165-0-1 (180-28384-3).

PCB-1016 failed the recovery criteria for the MS/MSD of sample SD-185-0-1 (180-28384-23). The presence of the '4' qualifier in the data indicates analytes where the concentration in the unspiked sample exceeded four times the spiking amount.

### **GENERAL CHEMISTRY**

No difficulties were encountered during the analysis.

### **SAMPLE SUMMARY**

Client: Tetra Tech, Inc. Job Number: 180-28384-1

			Date/Time	Date/Time
Lab Sample ID	Client Sample ID	Client Matrix	Sampled	Received
180-28384-1	SD-163-0-1	Sediment	12/18/2013 1155	12/23/2013 0900
180-28384-2	SD-164-0-1	Sediment	12/18/2013 1210	12/23/2013 0900
180-28384-3	SD-165-0-1	Sediment	12/18/2013 1235	12/23/2013 0900
180-28384-4	SD-166-0-1	Sediment	12/18/2013 1305	12/23/2013 0900
180-28384-5	SD-167-0-1	Sediment	12/18/2013 1510	12/23/2013 0900
180-28384-6	SD-168-0-1	Sediment	12/18/2013 1545	12/23/2013 0900
180-28384-7	SD-169-0-1	Sediment	12/18/2013 1610	12/23/2013 0900
180-28384-8	SD-170-0-1	Sediment	12/20/2013 1515	12/23/2013 0900
180-28384-9	SD-171-0-1	Sediment	12/19/2013 1045	12/23/2013 0900
180-28384-10	SD-172-0-3	Sediment	12/19/2013 1110	12/23/2013 0900
180-28384-11	SD-173-0-3	Sediment	12/19/2013 1130	12/23/2013 0900
180-28384-12	SD-174-0-1	Sediment	12/19/2013 1340	12/23/2013 0900
180-28384-13	SD-175-0-1	Sediment	12/19/2013 1430	12/23/2013 0900
180-28384-14	SD-176-0-3	Sediment	12/19/2013 1505	12/23/2013 0900
180-28384-15	SD-177-0-2	Sediment	12/19/2013 1540	12/23/2013 0900
180-28384-16	SD-178-0-1	Sediment	12/19/2013 1600	12/23/2013 0900
180-28384-17	SD-179-0-1	Sediment	12/19/2013 1645	12/23/2013 0900
180-28384-18	SD-180-0-1	Sediment	12/19/2013 1625	12/23/2013 0900
180-28384-19	SD-181-0-1	Sediment	12/20/2013 1000	12/23/2013 0900
180-28384-20	SD-182-0-3	Sediment	12/20/2013 1020	12/23/2013 0900
180-28384-21	SD-183-0-1	Sediment	12/20/2013 1040	12/23/2013 0900
180-28384-22	SD-184-0-2	Sediment	12/20/2013 1130	12/23/2013 0900
180-28384-23	SD-185-0-1	Sediment	12/20/2013 1200	12/23/2013 0900
180-28384-24	SD-186-0-1	Sediment	12/20/2013 1220	12/23/2013 0900
180-28384-25	SD-187-0-1	Sediment	12/20/2013 1245	12/23/2013 0900
180-28384-26	SD-188-0-1	Sediment	12/20/2013 1257	12/23/2013 0900
180-28384-27	SD-189-0-1	Sediment	12/20/2013 1250	12/23/2013 0900
180-28384-28	SD-190-0-1	Sediment	12/20/2013 1445	12/23/2013 0900
180-28384-29	SD-191-0-1	Sediment	12/20/2013 1450	12/23/2013 0900
180-28384-30	SD-192-0-1	Sediment	12/20/2013 1600	12/23/2013 0900
180-28384-31	SD-193-0-1	Sediment	12/20/2013 1610	12/23/2013 0900
180-28384-32	SD-194-0-1	Sediment	12/20/2013 1615	12/23/2013 0900

### FORM II GC/MS SEMI VOA SURROGATE RECOVERY

Lab Name:	TestAmerica Pittsburgh	Job No.:	180-28384-1
SDG No.:			
		_	

Matrix: Sediment Level: Low

GC Column (1): Rxi-5SilMS ID: 0.32 (mm)

			T			
Client Sample ID	Lab Sample ID	NBZ #	FBP	#	TPH	#
SD-163-0-1	180-28384-1	64 I	70	D	72	1
SD-164-0-1	180-28384-2	74 I	83	D	63	E
SD-165-0-1	180-28384-3	81 I	80	- D	67	Γ
SD-166-0-1	180-28384-4	0 X I	0 0	X D	0	ΧE
SD-167-0-1	180-28384-5	0 X I	0 0	X D	0	ΧГ
SD-168-0-1	180-28384-6	0 X I	0 0	X D	0	ΧГ
SD-169-0-1	180-28384-7	68 I	51	D	55	E
SD-170-0-1	180-28384-8	65 I	67	D	57	D
SD-171-0-1	180-28384-9	54	58		53	
SD-172-0-3	180-28384-10	54 I	53	D	47	D
SD-173-0-3	180-28384-11	65 I	67	D	59	D
SD-174-0-1	180-28384-12	65 I	69	D	58	Ε
SD-175-0-1	180-28384-13	63 I	62	D	64	Ι
SD-176-0-3	180-28384-14	57 I	57	D	45	D
SD-177-0-2	180-28384-15	40 I	42	D	39	Ī
SD-178-0-1	180-28384-16	53 I	56	D	52	D
SD-179-0-1	180-28384-17	63	61		46	
SD-180-0-1	180-28384-18	0 X I	0	X D	0	ΧĪ
SD-181-0-1	180-28384-19	0 X I	0	X D	0	ΧI
SD-182-0-3	180-28384-20	0. X I	0	X D	0	ΧI
SD-183-0-1	180-28384-21	51 I	61	D	54	Γ
SD-184-0-2	180-28384-22	62 I	70	D	61	I
SD-185-0-1	180-28384-23	62 I	74	D	61	Ι
SD-186-0-1	180-28384-24	76 I	90	D	76	Γ
SD-187-0-1	180-28384-25	77 I	92	D	83	Γ
SD-188-0-1	180-28384-26	84 I	98	D	84	D
SD-189-0-1	180-28384-27	74 I	94	D	79	
SD-190-0-1	180-28384-28	70 I	84	D <sub>:</sub>	89	Ι
SD-190-0-1 DL	180-28384-28 DL	62 I	83	D-	87	Ī
SD-191-0-1	180-28384-29	60 I	73	D	66	D
SD-191-0-1 DL	180-28384-29 DL	0 X I	0	X D		X D
SD-192-0-1	180-28384-30	52 I	67	D	56	Ė
SD-193-0-1	180-28384-31	58 I	72	D	89	Ε
SD-194-0-1	180-28384-32	65 I	80		66	Í
	MB 180-93179/1-A	75	68		74	

		QC LIMITS
NBZ =	Nitrobenzene-d5 (Surr)	27-110
FBP =	2-Fluorobiphenyl	28-108
TPH =	Terphenyl-d14 (Surr)	21-130

<sup>#</sup> Column to be used to flag recovery values

### FORM II GC/MS SEMI VOA SURROGATE RECOVERY

Lab Nam	e: <u>Te</u> stAmerica	Pittsburgh	Job No.	180-28384-1
SDG No.	:			
Matrix:	Sediment		Level:	Low

GC Column (1): Rxi-5SilMS ID: 0.32 (mm)

Client Sample ID	Lab Sample ID	NBZ	#	FBP	#	TPH	#
	MB 180-93531/1-A	57		59	+	55	
	LCS 180-93179/2-A	68	i	66		125	
	LCS 180-93531/2-A	57		60	-	58	
SD-163-0-1 MS	180-28384-1 MS	69	D	79	D	67	D
SD-185-0-1 MS	180-28384-23 MS	83	D	103	D	91	_ p
SD-163-0-1 MSD	180-28384-1 MSD	67	Ď	72	D	60	D
SD-185-0-1 MSD	180-28384-23 MSD	73	D	84	D	72	D

# Column to be used to flag recovery values

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: CCVIS 180-93752/22 Date Analyzed: 01/01/2014 04:44

Instrument ID: 71

GC Column: Rxi-5SilMS ID: 0.32(mm)

Lab File ID (Standard): S01010C2.D Heated Purge: (Y/N) N

Calibration ID: 12755

		DCB		NPT		ANT	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
12/24 HOUR STD		81807	6.23	381641	7.44	265775	9.07
UPPER LIMIT	·	163614	6.73	763282	7.94	531550	9.57
LOWER LIMIT		40904	5.73	190821	6.94	132888	8.57
LAB SAMPLE ID	CLIENT SAMPLE ID		•				
MB 180-93531/1-A		100562	6.22	459676	7.44	337997	9.06
LCS 180-93531/2-A		94791	6.23	432186	7.45	321735	9.06
180-28384-23 MS	SD-185-0-1 MS	102496	6.23	469936	7.45	334261	9.07
180-28384-23 MSD	SD-185-0-1 MSD	102718	6.23	467749	7.45	340218	9.07
180-28384-21	SD-183-0-1	84459	6.24	380362	7.45	269480	9.07
180-28384-22	SD-184-0-2	98329	6.23	448717	7.45	320106	9.06
180-28384-23	SD-185-0-1	92875	6.23	425521	7.45	308145	9.06
180-28384-24	SD-186-0-1	94962	6.23	436232	7.45	313779	9.06
180-28384-25	SD-187-0-1	99893	6.24	458246	7.44	330863	9.07
180-28384-26	SD-188-0-1	96925	6.24	440934	7.44	314348	9.06
180-28384-27	SD-189-0-1	98560	6.23	449313	7.44	318063	9.06
180-28384-28	SD-190-0-1	98108	6.24	448537	7.44	319559	9.06
180-28384-29	SD-191-0-1	97820	6,23	448311	7.45	317588	9.07
180-28384-30	SD-192-0-1	96741	6.24	445073	7.44	322461	9.07
180-28384-31	SD-193-0-1	103212	6.23	475585	7.44	339053	9.06
180-28384-32	SD-194-0-1	99925	6.24	451759	7.45	323303	9.07

DCB = 1,4-Dichlorobenzene-d4

NPT = Naphthalene-d8

ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

 $\ensuremath{\text{\#}}$  Column used to flag values outside QC limits

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: CCVIS 180-93752/22 Date Analyzed: 01/01/2014 04:44

Instrument ID: 71 GC Column: Rxi-5SilMS ID: 0.32(mm)

Lab File ID (Standard): S01010C2.D Heated Purge: (Y/N) N

Calibration ID: 12755

		PHN		CRY		PRY	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
12/24 HOUR STD		529007	10.43	650022	13.90	663127	16.79
UPPER LIMIT		1058014	10.93	1300044	14.40	1326254	17.29
LOWER LIMIT	m :	264504	9.93	325011	13.40	331564	16.29
LAB SAMPLE ID	CLIENT SAMPLE ID						
MB 180-93531/1-A		691656	10.43	841357	13.88	873602	16.78
LCS 180-93531/2-A		648601	10.43	784936	13.89	800504	16.78
180-28384-23 MS	SD-185-0-1 MS	693615	10.43	875273	13.89	969963	16.78
180-28384-23 MSD	SD-185-0-1 MSD	688931	10.43	896059	13.89	1021525	16.79
180-28384-21	SD-183-0-1	548984	10.43	682770	13.90	823897	16.80
180-28384-22	SD-184-0-2	630350	10.43	793827	13.90	953905	16.80
180-28384-23	SD-185-0-1	627324	10.43	824538	13.89	1001044	16.79
180-28384-24	SD-186-0-1	640443	10.43	848478	13.89	1037304	16.78
180-28384-25	SD-187-0-1	673158	10.43	895709	13.89	1109145	16.79
180-28384-26	SD-188-0-1	642180	10.43	864756	13.89	1098047	16.79
180-28384-27	SD-189-0-I	655592	10.43	879684	13.89	1136630	16.79
180-28384-28	SD-190-0-1	644328	10.43	817781	13.90	1173363	16.80
180-28384-29	SD-191-0-1	634911	10.43	821271	13.91	1195692	16.82
180-28384-30	SD-192-0-1	655710	10.43	883014	13.89	1175551	16.80
180-28384-31	SD-193-0-1	704065	10.43	940080	13.89	1246020	16.79
180-28384-32	SD-194-0-1	668717	10.43	913984	13.90	1216412	16.80

PHN = Phenanthrene-d10

CRY = Chrysene-d12

PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: CCVIS 180-93785/3 Date Analyzed: 01/04/2014 07:53

Instrument ID: 71 GC Column: Rxi-5SilMS ID: 0.32(mm)

Lab File ID (Standard): S01040C1.D Heated Purge: (Y/N) N

Calibration ID: 12755

	DCB		NPT		ANT	
	AREA #	RT #	AREA #	RT #	AREA #	RT #
12/24 HOUR STD	77376	6,17	311581	7.39	194573	9.01
UPPER LIMIT	154752	6.67	623162	7.89	389146	9.51
LOWER LIMIT	38688	5.67	155791	6.89	97287	8.51
LAB SAMPLE ID CLIENT SAMPLE ID						
180-28384-28 DL SD-190-0-1 DL	67214	6.18	272161	7.39	182839	9.01

DCB = 1,4-Dichlorobenzene-d4

NPT = Naphthalene-d8

ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

Lab Name:	TestAmerica Pittsburgh	Job No.: 180-28384-1				
SDG No.:						
Sample No.	: CCVIS 180-93785/3	Date Analyzed: 01/04/2014 07:53				
Instrument	: ID: <u>71</u>	GC Column: Rxi-5SilMS ID:	0.32 (mm)			
Lab File 1	ID (Standard): S01040C1.D	Heated Purge: (Y/N) N				
Calibratio	on ID: 12755					

	PHN		CRY		PRY		
	AREA #	RT #	AREA #	RT #	AREA #	RT	
12/24 HOUR STD	377760	10.38	426352	13.82	379629	16.69	
UPPER LIMIT	755520	10.88	852704	14.32	759258	17.19	
LOWER LIMIT	188880	9.88	213176	13.32	189815	16.19	

364362

10.38

455661

13.81

497985

16.68

PHN = Phenanthrene-d10

CRY = Chrysene-d12

180-28384-28 DL

PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

SD-190-0-1 DL

 $\ensuremath{\sharp}$  Column used to flag values outside QC limits

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: CCVIS 180-93945/2 Date Analyzed: 01/08/2014 01:27

Instrument ID: 722 GC Column:

GC Column: Rxi-5SilMS ID: 0.32(mm)

Lab File ID (Standard): F01080C1.D Heated Purge: (Y/N) N

Calibration ID: 13164

			DCB		NPT		ANT		
			AREA #	RT #	AREA #	RT #	AREA #	RT #	
12/24 HOUR STD			54457	6.15	234208	7.41	139630	9.11	
UPPER LIMIT		_ i	108914	6.65	468416	7.91	279260	9.61	
LOWER LIMIT			27229	5.65	117104	6.91	69815	8.61	
LAB SAMPLE ID	CLIENT SAMPLE ID	=		1					
180-28384-29 DL	SD-191-0-1 DL		53786	6.15	232986	7.42	149598	9.11	

DCB = 1,4-Dichlorobenzene-d4

NPT = Naphthalene-d8

ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

### 

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: CCVIS 180-93945/2 Date Analyzed: 01/08/2014 01:27

Instrument ID: 722 GC Column: <u>Rxi-5SilMS</u> ID: 0.32(mm)

Lab File ID (Standard): F01080C1.D Heated Purge: (Y/N) N

Calibration ID: 13164

	PHN		CRY		PRY		
	AREA #	RT #	AREA #	RT #	AREA #	RT #	
12/24 HOUR STD	256701	10.55	276001	14.39	237474	17.55	
UPPER LIMIT	513402	11.05	552002	14.89	474948	18.05	
LOWER LIMIT	128351	10.05	138001	13.89	118737	17.05	
LAB SAMPLE ID CLIENT SAMPLE ID							
180-28384-29 DL SD-191-0-1 DL	263447	10.55	292098	14.36	285280	17.50	

PHN = Phenanthrene-d10

CRY = Chrysene-d12

PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.: Sample No.: CCVIS 180-93613/22 Date Analyzed: 12/31/2013 10:31

Instrument ID: 732 GC Column: Rxi-5SilMS ID: 0.32(mm)

Lab File ID (Standard): D12310CC.D Heated Purge: (Y/N) N

Calibration ID: 12823

		DCB		NPT		ANT	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
12/24 HOUR STD		146077	6.15	593830	7.44	44 373677	
UPPER LIMIT		292154	6.65	1187660	7.94	747354	9.65
LOWER LIMIT		73039	5.65	296915	6.94	186839	8.65
LAB SAMPLE ID	CLIENT SAMPLE ID						
MB 180-93179/1-A		154463	6.13	625757	7.42	395469	9.13
LCS 180-93179/2-A		152717	6.12	613059	7.41	399109	9.13
180-28384-1 MS	SD-163-0-1 MS	157285	6.12	629331	7.42	408727	9.13
180-28384-1 MSD	SD-163-0-1 MSD	160765	6.13	635904	7.42	410679	9.13
180-28384-1	SD-163-0-1	172869	6.13	707409	7.42	439215	9.13
180-28384-2	SD-164-0-1	167636	6.12	692592	7.42	431961	9.13
180-28384-3	SD-165-0-1	159091	6.13	635444	7.42	419498	9.13
180-28384-4	SD-166-0-1	162911	6.13	665898	7.42	446513	9.14
180-28384-5	SD-167-0-1	162581	6.12	669832	7.42	436395	9.13
180-28384-6	SD-168-0-1	150752	6.12	617067	7.41	405845	9.13
180-28384-7	SD-169-0-1	197833	6.12	831934	7.42	540243	9.13
180-28384-8	SD-170-0-1	198383	6.12	799184	7.41	521027	9.13
180-28384-9	SD-171-0-1	204156	6.13	831963	7.42	528125	9.13
180-28384-10	SD-172-0-3	194005	6.12	813387	7.42	518524	9.13
180-28384-11	SD-173-0-3	197048	6.12	823834	7.42	543624	9.13
180-28384-12	SD-174-0-1	189846	6.12	783286	7.42	519442	9.13
180-28384-13	SD-175-0-1	188798	6.12	762865	7.41	503822	9.13
180-28384-14	SD-176-0-3	190445	6.12	786927	7.41	501396	9.13
180-28384-15	SD-177-0-2	207954	6.13	832407	7.42	545365	9.13
180-28384-16	SD-178-0-1	208866	6.12	833391	7.42	544653	9.13
180-28384-17	SD-179-0-1	190625	6.12	787227	7.42	506133	9.13

DCB = 1,4-Dichlorobenzene-d4

NPT = Naphthalene-d8

ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

 $\mbox{\#}$  Column used to flag values outside QC limits

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: CCVIS 180-93613/22 Date Analyzed: 12/31/2013 10:31

Instrument ID: 732 GC Column: Rxi-5SilMS ID: 0.32(mm)

Lab File ID (Standard): D12310CC.D Heated Purge: (Y/N) N

Calibration ID: 12823

		PHN		CRY	. ]	PRY	
		AREA #	RT #	AREA #	RT #	AREA #	RT ‡
12/24 HOUR STD		677230	10.58	792212	14.32	663428	17.21
UPPER LIMIT		1354460	11.08	1584424	14.82	1326856	17.71
LOWER LIMIT		338615	10.08	396106	13.82	331714	16.71
LAB SAMPLE ID	CLIENT SAMPLE ID		!				
MB 180-93179/1-A		727395	10.57	855339	14.29	763573	17.18
LCS 180-93179/2-A		757536	10.56	892890	14.29	751960	17.17
180-28384-1 MS	SD-163-0-1 MS	771814	10.57	928598	14.30	876378	17.18
180-28384-1 MSD	SD-163-0-1 MSD	777453	10.57	1010165	14.30	974595	17.18
180-28384-1	SD-163-0-1	815893	10.57	1028757	14.31	1039007	17.19
180-28384-2	SD-164-0-1	782158	10.57	1010873	14.29	1035145	17.17
180-28384-3	SD-165-0-1	806201	10.57	971798	14.30	967067	17.19
180-28384-4	SD-166-0-1	821512	10.57	1016240	14.31	1048144	17.18
180-28384-5	SD-167-0-1	801581	10.57	1053680	14.30	1089120	17.18
180-28384-6	SD-168-0-1	757668	10.56	983735	14.29	1040244	17.17
180-28384-7	SD-169-0-1	990004	10.57	1330221	14.29		17.18
180-28384-8	SD-170-0-1	974947	10.56	1241347	14.28	1434416*	17.16
180-28384-9	SD-171-0-1	1007574	10.57	1305918	14.31	1545689*	17.19
180-28384-10	SD-172-0-3	993257	10.57	1234921	14.29	1496516*	17.18
180-28384-11	SD-173-0-3	1011966	10.57	1311268	14.30	1588006*	17.19
180-28384-12	SD-174-0-1	989576	10.57	1273244	14.31	1587998*	17.19
180-28384-13	SD-175-0-1	927398	10.57	1245091	14.30	1532160*	17.18
180-28384-14	SD-176-0-3	969462	10.56	1342030	14.29	1635796*	17.17
180-28384-15	SD-177-0-2	1033002	10.57	1409372	14.31	1750837	17.20
180-28384-16	SD-178-0-1	1058398	10.57	1439958	14.30	1750128*	17.18
180-28384-17	SD-179-0-1	977214	10.57	1402126	14.30	1688397*	17.19

PHN = Phenanthrene-d10

CRY = Chrysene-d12

PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

 $\ensuremath{\mathtt{\#}}$  Column used to flag values outside QC limits

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: CCVIS 180-93710/10 Date Analyzed: 01/02/2014 12:15

Instrument ID: 732 GC Cc

GC Column: Rxi-5SilMS ID: 0.32(mm)

Lab File ID (Standard): D01020CC.D Heated Purge: (Y/N) N

Calibration ID: 12823

		DCB		NPT		ANT	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
12/24 HOUR STD		187827	6.13	691397	7.43	396729	9.16
UPPER LIMIT		375654	6.63	1382794	7.93	793458	9.66
LOWER LIMIT	93914	5.63	345699	6.93	198365	8.66	
LAB SAMPLE ID	CLIENT SAMPLE ID				i i		
180-28384-18	SD-180-0-1	213650	6.12	789029	7.42	437781	9.16
180-28384-19	SD-181-0-1	229336	6.12	886719	7.43	501642	9.16
180-28384-20	SD-182-0-3	226825	6.12	845333	7.43	470738	

DCB = 1,4-Dichlorobenzene-d4

NPT = Naphthalene-d8

ANT = Acenaphthene-d10

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: CCVIS 180-93710/10 Date Analyzed: 01/02/2014 12:15

Instrument ID: 732 GC Column: Rxi-5SilMS ID: 0.32(mm)

Lab File ID (Standard): D01020CC.D Heated Purge: (Y/N) N

Calibration ID: 12823

		PHN		CRY		PRY		
		AREA #	RT #	AREA #	RT #	AREA #	RT #	
12/24 HOUR STD		659904	10.61	664044	14.39	534460	17.29	
UPPER LIMIT		1319808	11.11	1328088	14.89	1068920	17.79	
LOWER LIMIT		329952	10.11	332022	13.89	267230	16.79	
LAB SAMPLE ID	CLIENT SAMPLE ID							
180-28384-18	SD-180-0-1	667938	10.61	626406	14.41	576992	17.32	
180-28384-19	SD-181-0-1	758727	10.61	661054	14.41	645429	17.32	
180-28384-20	SD-182-0-3	709972	10.61	667487	14.41	679232	17.32	

PHN = Phenanthrene-d10

CRY = Chrysene-d12

PRY = Perylene-d12

Area Limit = 50%-200% of internal standard area RT Limit =  $\pm$  0.5 minutes of internal standard RT

# Column used to flag values outside QC limits

### FORM IV GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Pittsburgh	Job No.: 180-28384-1
SDG No.:	
Lab File ID: D1231003.D	Lab Sample ID: MB 180-93179/1-A
Matrix: Sediment	Date Extracted: 12/24/2013 03:15
Instrument ID: 732	Date Analyzed: 12/31/2013 10:59
Level: (Low/Med) Low	

### THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

		LAB	
CLIENT SAMPLE ID	LAB SAMPLE ID	FILE ID	DATE ANALYZED
	LCS 180-93179/2-A	D1231004.D	12/31/2013 11:51
SD-163-0-1 MS	180-28384-1 MS	D1231005.D	12/31/2013 14:03
SD-163-0-1 MSD	180-28384-1 MSD	D1231006.D	12/31/2013 14:29
SD-163-0-1	180-28384-1	D1231007.D	12/31/2013 14:56
SD-164-0-1	180-28384-2	D1231008.D	12/31/2013 15:22
SD-165-0-1	180-28384-3	D1231009.D	12/31/2013 15:49
SD-166-0-1	180-28384-4	D1231010.D	12/31/2013 16:15
SD-167-0-1	180-28384-5	D1231011.D	12/31/2013 16:41
SD-168-0-1	180-28384-6	D1231012.D	12/31/2013 17:08
SD-169-0-1	180-28384-7	D1231013.D	12/31/2013 17:34
SD-170-0-1	180-28384-8	D1231014.D	12/31/2013 18:00
SD-171-0-1	180-28384-9	D1231015.D	12/31/2013 18:26
SD-172-0-3	180-28384-10	D1231016.D	12/31/2013 18:53
SD-173-0-3	180-28384-11	D1231017.D	12/31/2013 19:19
SD-174-0-1	180-28384-12	D1231018.D	12/31/2013 19:46
SD-175-0-1	180-28384-13	D1231019.D	12/31/2013 20:12
SD-176-0-3	180-28384-14	D1231020.D	12/31/2013 20:38
SD-177-0-2	180-28384-15	D1231021.D	12/31/2013 21:04
SD-178-0-1	180-28384-16	D1231022.D	12/31/2013 21:31
SD-179-0-1	180-28384-17	D1231023.D	12/31/2013 21:57
SD-180-0-1	180-28384-18	D0102001.D	01/02/2014 19:05
SD-181-0-1	180-28384-19	D0102002.D	01/02/2014 19:32
SD-182-0-3	180-28384-20	D0102003.D	01/02/2014 19:58

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SDG No.:					***************************************
Batch Number:	93179	Batch Start Date: 12/24/13 03:15		Batch Analyst:	Batch Analyst: Geehring, Kevin
Batch Method:	3541	Batch End Date: 12/24/13 07:48	/24/13 07:48		

OPQL8270SURi 00012	20 uL	20 nT	20 nF	20 nF	50 uL	20 nr	50 uL	50 uL	50 uL	50 uL	50 uL	50 uL	20 uL	50 uL	50 uL	50 uL					
OPLVISPKMIX1i 00021		50 uL	50 uL	50 uL																	
InitialAmount	30.0 g	30.0 g	30.0 g	30.1 g	30.1 g	30.0 g	30.1 g	30.0 g	30.0 g	30.1 g	30.1 g	30.0	30.0 g	30.1 g	30.0 g	30.0 g	30.0 g	30.1 g	30.2 g	30.0 g	30.0 g
FinalAmount	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	1.0 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL
Basis			E	EH		F	E	T.	L.	E	EH	E	E-I	E	L	£ .	E+	E-	L	<u></u>	E
Method Chain	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D	3541, 8270D LL
Client Sample ID			SD-163-0-1	SD-163-0-1	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-167-0-1	SD-168-0-1	SD-169-0-1	SD-170-0-1	SD-171-0-1	SD-172-0-3	SD-173-0-3	SD-174-0-1	SD-175-0-1	SD-176-0-3	SD-177-0-2	SD-178-0-1	SD-179-0-1
Lab Sample ID	MB 180-93179/1	LCS 180-93179/2	180-28384-A-1	180-28384-A-1	180-28384-A-1	180-28384-A-2	180-28384-A-3	180-28384-A-4	180-28384-A-5	180-28384-A-6	180-28384-A-7	180-28384-A-8	180-28384-A-9	180-28384-A-10	180-28384-A-11	180-28384-A-12	180-28384-A-13	180-28384-A-14	180-28384-A-15	180-28384-A-16	180-28384-A-17

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.

のいいが、神神の歌曲、空間のはなりのでは、いちにはなって、この歌をはながら、の では語いない しになる こじゅう かん 最後間でき

 $\sim$ 

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# GC/MS SEMI VOA BATCH WORKSHEET

Geehring, Kevin Batch Analyst: OPQL8270SURi 00012 50 uL 50 uL 50 uL OPLVISPKMIX11 03:15 12/24/13 07:48 InitialAmount 12/24/13 30.1 g 30.0 g 30.1 g Job No.: 180-28384-1 Batch Start Date: Batch End Date: Methylene chloride FinalAmount 0.5 mL 0.5 mL Mec12 / Acetone 0.5 mL 32 Degrees 1120122641 1053215 1047333 1055012 769170 SOX Batch Notes kg Basis kg Method Chain 3541, 8270D LL 3541, 8270D LL 3541, 8270D LL TestAmerica Pittsburgh Person's name who did the concentration Client Sample ID Basis Description Person's name who did the prep Uncorrected N-evap Temperature SD-180-0-1 SD-181-0-1 SD-182-0-3 93179 3541 Exchange Solvent Lot # Magnesium Sulfate Lot Exchange Solvent Name Na2SO4 Lot Number Batch Method: Batch Number: T | Total/NA Lab Sample ID 180-28384-A-18 180-28384-A-20 180-28384-A-19 Batch Comment Lab Name: Solvent Lot Balance ID SDG No.: N-evap # Solvent Basis

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.

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### FORM I GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1 SDG No.: Lab Sample ID: MB 180-93179/1-A Client Sample ID: Matrix: Sediment Lab File ID: D1231003.D Analysis Method: 8270D LL Date Collected: Extract. Method: 3541 Date Extracted: 12/24/2013 03:15 Sample wt/vol: 30.0(g) Date Analyzed: 12/31/2013 10:59 Dilution Factor: 1 Con. Extract Vol.: 0.5(mL) Injection Volume: 2(uL) Level: (low/med) Low % Moisture: GPC Cleanup: (Y/N) N Analysis Batch No.: 93613 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT Q	RL	MDL
120-12-7	Anthracene	ND	3.4	0.33
56-55-3	Benzo[a]anthracene	ND	3.4	0.42
205-99-2	Benzo[b] fluoranthene	ND	3.4	0.52
207-08-9	Benzo[k]fluoranthene	ND	3.4	0.67
191-24-2	Benzo[g,h,i]perylene	ND	3.4	0.33
50-32-8	Benzo[a]pyrene	ND	3.4	0.33
218-01-9	Chrysene	ND	3.4	0.40
53-70-3	Dibenz(a,h)anthracene	ND	3.4	0.37
206-44-0 .	Fluoranthene	ND	3.4	0.36
86-73-7	Fluorene	ND	3.4	0.44
193-39-5	Indeno[1,2,3-cd]pyrene	ND	3.4	0.34
85-01-8	Phenanthrene	ND	3.4	0.53
129-00-0	Pyrene	ND	3.4	0.34
83-32-9	Acenaphthene	ND	3.4	0.32
208-96-8	Acenaphthylene	ND	3.4	0.38
91-20-3	Naphthalene	ND	3.4	0.29
91-57-6	2-Methylnaphthalene	ND	3.4	0.30

CAS NO.	SURROGATE	%REC Q	LIMITS
4165-60-0	Nitrobenzene-d5 (Surr)	75	27-110
321-60-8	2-Fluorobiphenyl	68	28-108
1718-51-0	Terphenyl-d14 (Surr)	74	21-130

### FORM III GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.: \_\_\_\_

Matrix: Sediment Level: Low Lab File ID: D1231004.D

Lab ID: LCS 180-93179/2-A Client ID:

	SPIKE	LCS	LCS	QC	
	ADDED	CONCENTRATION	90	LIMITS	#
COMPOUND	(ug/Kg)	(ug/Kg)	REC	REC	
Anthracene	333	221	66	43-111	
Benzo[a]anthracene	333	235	70	45-110	
Benzo[b] fluoranthene	333	233	70	37-108	
Benzo[k] fluoranthene	333	227	68	39-115	
Benzo[g,h,i]perylene	333	249	75	35-127	
Benzo[a]pyrene	333	247	74	42-114	
Chrysene	333	239	72	44-108	
Dibenz(a,h)anthracene	333	249	75	34-131	
Fluoranthene	333	228	69	40-118	
Fluorene	333	221	66	43-110	
<pre>Indeno[1,2,3-cd]pyrene</pre>	333	241	72	34-130	
Phenanthrene	333	216	65	41-107	
Pyrene	333	235	71	39-113	
Acenaphthene	333	. 216	65	42-104	
Acenaphthylene	333	226	68	43-117	
Naphthalene	333	200	60	42-104	
2-Methylnaphthalene	333	201	60	43-105	

<sup>#</sup> Column to be used to flag recovery and RPD values FORM III 8270D LL

### FORM III GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.: \_\_\_\_

Matrix: Sediment Level: Low Lab File ID: S0101003.D

Lab ID: 180-28384-23 MS Client ID: SD-185-0-1 MS

	SPIKE	SAMPLE	MS	MS	QC :	
	ADDED	CONCENTRATION	CONCENTRATION	용	LIMITS	#
COMPOUND	(ug/Kg)	(ug/Kg)	(ug/Kg)	REC	REC	•
Anthracene	445	230	818	131	43-111	F1
Benzo[a]anthracene	445	1300	1400	21	45-110	F1
Benzo[b] fluoranthene	445	1500	1510	-0.4	37-108	F1
Benzo[k]fluoranthene	445	570	939	82	39-115	
Benzo[g,h,i]perylene	445	1100	1400	59	35-127	
Benzo[a]pyrene	445	1200	1360	29	42-114	F1
Chrysene	445	1600	1810	47	44-108	
Dibenz(a,h)anthracene	445	270	706	98	34-131	
Fluoranthene	445	3600	3350	-58	40-118	(4)
Fluorene	445	87 J	671	(131	43-110	F1
<pre>Indeno[1,2,3-cd]pyrene</pre>	445	1000	1190	40	34-130	
Phenanthrene	445	1100	1860	181	41-107	F1
Pyrene	445	1900	2000	- 20	39-113	(4)
Acenaphthene	445	53 J	582	119	42-104	F1
Acenaphthylene	445	94 Ј	536	99	43-117	
Naphthalene	445	43 J	464	94	42-104	
2-Methylnaphthalene	445	21 J	517	111	43-105	F1

<sup>#</sup> Column to be used to flag recovery and RPD values FORM III 8270D LL

### FORM III GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name	e: TestAmerica Pitt	sburgh	Job No.: 180-28384-1					
SDG No.:				_				
Matrix:	Sediment	Level: Low	Lab File ID: S0101004.D					
Lab ID:	180-28384-23 MSD		Client ID: SD-185-0-1 MSD					

	SPIKE	MSD	MSD		QC LI	MITS
	ADDED	CONCENTRATION	8	8		#
COMPOUND	(ug/Kg)	(ug/Kg)	REC	RPD	RPD_	REC
Anthracene	447	497	59	(49)	35	43-111 F2
Benzo[a]anthracene	447	916	(-89		(313)	45-110 F1 F2
Benzo[b]fluoranthene	447	1070	<u>-98</u>	(34)	<b>2</b>	37-108 F1 F2
Benzo[k]fluoranthene	447	689	26		42	39-115 F1
Benzo[g,h,i]perylene	447	1050	- <del>- 2</del> 0	29	21	35-127 F1 F2
Benzo[a]pyrene	447	903	74		31	42-114 F1 F2
Chrysene	447	1130	€105	<b>(46)</b>	31	44-108 F1 F2
Dibenz(a,h)anthracene	447	563	<u>_65</u>	23	32	34-131
Fluoranthene	447	1960	<del>-369</del>	ノ (52)	23	40-118 <b>/</b> 4 <b>)</b> F2
Fluorene	447	474	87	34	37	43-110
Indeno[1,2,3-cd]pyrene	447	923	(-20	<b>)</b> 25	30	34-130 F1
Phenanthrene	447	848	<u>-46</u>		20	41-107 F1 F2
Pyrene	447	1130 -	<b>(</b> -170	) (56)	28	39-113 <b>(4)</b> F2
Acenaphthene	447	403	78	- B	34	42-104 F2
Acenaphthylene	447	431	75		36	43-117
Naphthalene	447	371	73	22	25	42-104
2-Methylnaphthalene	447	412	87	23	34	43-105

 $<sup>\</sup>mbox{\#}$  Column to be used to flag recovery and RPD values FORM III 8270D  $\mbox{LL}$ 

### FORM IV GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab File ID: S0101001.D Lab Sample ID: MB 180-93531/1-A

Matrix: Sediment Date Extracted: 12/31/2013 03:40

Instrument ID: 71 Date Analyzed: 01/01/2014 05:41

Level: (Low/Med) Low

### THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
LCS 180-93531/2-A	S0101002.D	01/01/2014 06:35
180-28384-23 MS	S0101003.D	01/01/2014 08:22
180-28384-23 MSD	S0101004.D	01/01/2014 08:48
180-28384-21	S0101005.D	01/01/2014 09:15
180-28384-22	S0101006.D	01/01/2014 09:42
180-28384-23	S0101007.D	01/01/2014 10:08
180-28384-24	S0101008.D	01/01/2014 10:35
180-28384-25	S0101009.D	01/01/2014 11:01
180-28384-26	S0101010.D	01/01/2014 11:28
180-28384-27	S0101011.D	01/01/2014 11:55
180-28384-28	S0101012.D	01/01/2014 12:22
180-28384-29	S0101013.D	01/01/2014 12:48
180-28384-30	S0101014.D	01/01/2014 13:15
180-28384-31	S0101015.D	01/01/2014 13:42
180-28384-32	S0101016.D	01/01/2014 14:09
180-28384-28 DL	S0104012.D	01/04/2014 13:38
180-28384-29 DL	F0108001.D	01/08/2014 02:23
	LCS 180-93531/2-A  180-28384-23 MS  180-28384-21  180-28384-21  180-28384-22  180-28384-24  180-28384-25  180-28384-26  180-28384-27  180-28384-29  180-28384-30  180-28384-31  180-28384-32  180-28384-32  180-28384-38 DL	LCS 180-93531/2-A S0101002.D  180-28384-23 MS S0101003.D  180-28384-23 MSD S0101004.D  180-28384-21 S0101005.D  180-28384-22 S0101006.D  180-28384-23 S0101007.D  180-28384-24 S0101008.D  180-28384-25 S0101009.D  180-28384-26 S0101010.D  180-28384-27 S010101.D  180-28384-28 S0101012.D  180-28384-29 S0101013.D  180-28384-30 S0101015.D  180-28384-31 S0101015.D  180-28384-32 S0101016.D  180-28384-28 DL S0104012.D

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## GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: Tes	TestAmerica Pittsburgh	Job No.: 180-28384-1		
SDG No.:				
Batch Number:	93531	Batch Start Date: 12/31/13 03:40	Batch Analyst:	Batch Analyst: Geehring, Kevin
Batch Method:	3541	Batch End Date: 12/31/13 07:38		

						11 - 1 to the last annual management of the last of th										THE PROPERTY AND ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY AND ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY ADDRESS OF THE PROPERTY A
OPQL8270SURi 00012	50 uL	50 uL	50 uL	50 uL	50 uL	50 uL	20 uL	50 uL	50 uL	50 uL	50 uL	50 uL	50 uL	50 uL	50 uL	50 uL
OPLVISPKMIX1i 00021		20 uL	20 nT	50 uL												
InitialAmount	30.0 g	30.0 g	30.1 g	30.0 g	30.1 g	30.1 g	30.0 9	30.0 g	30.1 g	30.0 9	30.0 g	30.0 g	30.1 9	30.2 g	30.1 9	30.09
FinalAmount	0.5 mL	0.5 mL	1.0 mL	1.0 mL	0.5 mL	0.5 mL	1.0 mL	2.0 mL	3.0 mL	2.0 mL	2.0 mL	2.0 mL	0.5 mL	0.5 mL	0.5 mL	0.5 mL
Basis			L	H	T	T	L	£	E+	 E+	₽	H	 ⊢	F	F	H
Method Chain	3541, 8270D LL	3541, 8270D LL	3541, 8270D LL	3541, 8270D LL	3541, 8270D LL	3541, 8270D LL	3541, 8270D LL	3541, 8270D	3541, 8270D	3541, 8270D LL	3541, 8270D LL	3541, 8270D LL	3541, 8270D LL	3541, 8270D LL	3541, 8270D	3541, 8270D LL
Client Sample ID			SD-185-0-1	SD-185-0-1	SD-183-0-1	SD-184-0-2	SD-185-0-1	SD-186-0-1	SD-187-0-1	SD-188-0-1	SD-189-0-1	SD-190-0-1	SD-191-0-1	SD-192-0-1	SD-193-0-1	SD-194-0-1
Lab Sample ID	MB 180-93531/1	LCS 180-93531/2	180-28384-A-23 MS	180-28384-A-23 MSD	180-28384-A-21	180-28384-A-22	180-28384-A-23	180-28384-A-24	180-28384-A-25	180-28384-A-26	180-28384-A-27	180-28384-A-28	180-28384-A-29	180-28384-A-30	180-28384-A-31	180-28384-A-32

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 I GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Client Sample ID: Lab Sample ID: MB 180-93531/1-A

Matrix: Sediment Lab File ID: S0101001.D

Analysis Method: 8270D LL Date Collected:

Extract. Method: 3541 Date Extracted: 12/31/2013 03:40

Sample wt/vol: 30.0(g) Date Analyzed: 01/01/2014 05:41

Con. Extract Vol.: 0.5(mL) Dilution Factor: 1

Injection Volume: 2(uL) Level: (low/med) Low

% Moisture: GPC Cleanup:(Y/N) N

Analysis Batch No.: 93752 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT Q	RL	MDL
120-12-7	Anthracene	ND	3.4	0.33
56-55-3	Benzo[a]anthracene	ND	3.4	0.42
205-99-2	Benzo[b] fluoranthene	ND	3.4	0.52
207-08-9	Benzo[k]fluoranthene	ND	3.4	0.67
191-24-2	Benzo[g,h,i]perylene	ND	3.4	0.33
50-32-8	Benzo[a]pyrene	ΝĐ	3.4	0.33
218-01-9	Chrysene	. ND	3.4	0.40
53-70-3	Dibenz(a,h)anthracene	ND	3.4	0.37
206-44-0	Fluoranthene	ND	3.4	0.36
86-73-7	Fluorene	ND	3.4	0.44
193-39-5	Indeno[1,2,3-cd]pyrene	ND ·	3.4	0.34
85-01-8	Phenanthrene	ND	3.4	0.53
129-00-0	Pyrene	ND	3.4	0.34
83-32-9	Acenaphthene	ND	3.4	0.32
208-96-8	Acenaphthylene	ND	3.4	0.38
91-20-3	Naphthalene	ND	3.4	0.29
91-57-6	2-Methylnaphthalene	ND	3.4	0.30

CAS NO.	SURROGATE	%REC	Q	LIMITS
4165-60-0	Nitrobenzene-d5 (Surr)	57		27-110
321-60-8	2-Fluorobiphenyl	59		28-108
1718-51-0	Terphenyl-d14 (Surr)	55		21-130

### FORM III GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name	: TestAmerica Pitts	burgh	Job No.: 180	-28384-1
SDG No.:				·
Matrix:	Sediment	Level: Low	Lab File ID:	S0101002.D

Lab ID: LCS 180-93531/2-A Client ID: \_\_\_\_\_

	SPIKE	LCS	LCS	QC	
	ADDED	CONCENTRATION	8	LIMITS	#
COMPOUND	(ug/Kg)	(ug/Kg)	REC	REC	
Anthracene	333	232	70	43-111	
Benzo[a]anthracene	333	239	72	45-110	
Benzo[b] fluoranthene	333	217	65	37-108	
Benzo[k] fluoranthene	333	221	66	39-115	
Benzo[g,h,i]perylene	333	250	75	35-127	
Benzo[a]pyrene	333	234	70	42-114	
Chrysene	333	250	75	44-108	
Dibenz(a,h)anthracene	333	250	75	34-131	
Fluoranthene	333	271	81	40-118	
Fluorene	333	230	69	43-110	
<pre>Indeno[1,2,3-cd]pyrene</pre>	333	245	74	34-130	
Phenanthrene	333	219	66	41-107	
Pyrene	333	189	57	39-113	
Acenaphthene	333	208	62	42-104	
Acenaphthylene	333	218	65	43-117	
Naphthalene	333	209	63	42-104	
2-Methylnaphthalene	333	225	68	43-105	

 $<sup>\</sup>mbox{\#}$  Column to be used to flag recovery and RPD values FORM III 8270D LL

### FORM III GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Matrix: Sediment Level: Low Lab File ID: D1231005.D

Lab ID: 180-28384-1 MS Client ID: SD-163-0-1 MS

		т	T			
	SPIKE	SAMPLE	MS	MS	QC	
	ADDED	CONCENTRATION	CONCENTRATION	ક	LIMITS	#
COMPOUND	(ug/Kg)	(ug/Kg)	(ug/Kg)	REC !	REC	
Anthracene	434	97 J	1120	235	43-111	F1
Benzo[a]anthracene	434	530	2490	452	45-110	F1
Benzo[b] fluoranthene	434	700	2160	338	37-108	F1
Benzo[k]fluoranthene	434	260	1050	182	39-115	F1
Benzo[g,h,i]perylene	434	480	1930	335	35-127	F1
Benzo[a]pyrene	434	500	2110	370	42-114	F1
Chrysene	434	710	2850	494	44-108	F1
Dibenz(a,h)anthracene	434	110 J	617	118	34-131	
Fluoranthene	434	1400	5070	847	40-118	F1
F1uorene	434	52 J	894	194	43-110	F1
Indeno[1,2,3-cd]pyrene	434	390	1430	240	34-130	F1
Phenanthrene	434	990	6190	1197	41-107	F1
Pyrene	434	1100	6650	1268	39-113	F1
Acenaphthene	434	64 J	592	122	42-104	F1
Acenaphthylene	434	69 J	734	153	43-117	F1
Naphthalene	434	ND	468	108	42-104	F1
2-Methylnaphthalene	434	ND	592	136	43-105	F1

 $<sup>\</sup>mbox{\#}$  Column to be used to flag recovery and RPD values FORM III 8270D LL

### FORM III GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Matrix: Sediment Level: Low Lab File ID: D1231006.D

Lab ID: 180-28384-1 MSD Client ID: SD-163-0-1 MSD

	SPIKE	MSD CONCENTRATION	MSD %	90	QC LI	MITS	#
COMPOUND	(ug/Kg)	(ug/Kg)	REC	RPD	RPD :	REC	
Anthracene	433	1080	/22X	4	35	43-111	F1
Benzo[a]anthracene	433	2290	407	8	31	45-110	F1
Benzo[b] fluoranthene	433	2190	345	1	28	37-108	F1
Benzo[k]fluoranthene	433	805	126	26	42	39-115	F1
Benzo[g,h,i]perylene	433	1920	332	1	21	35-127	F1
Benzo[a]pyrene	433	2020	351	4	31	42-114	F1
Chrysene	433	2660	1451	7;	31	44-108	F1
Dibenz(a,h)anthracene	433	589	112	5	32	34-131	
Fluoranthene	433	4860	801	4	23	40-118	F1
Fluorene	433	894	194	0	37	43-110	F1
<pre>Indeno[1,2,3-cd]pyrene</pre>	433	1500	255	4	30	34-130	F1
Phenanthrene	433	5820	1117	6	20	41-107	F1
Pyrene	433	5910	1102	12	28	39-113	F1
Acenaphthene	433	593	122	0	34	42-104	F1
Acenaphthylene	433	744	156	1 1	36	43-117	F1
Naphthalene	433	431	99	6 V6	25	42-104	
2-Methylnaphthalene	433	563	130	5	34	43-105	F1

<sup>#</sup> Column to be used to flag recovery and RPD values FORM III 8270D  ${\rm LL}$ 

### FORM V GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab File ID: F1225DF1.D DFTPP Injection Date: 12/25/2013

Instrument ID: 722 DFTPP Injection Time: 19:08

Analysis Batch No.: 93544

M/E	ION ABUNDANCE CRITERIA		ATIVE DANCE
51	30.0 - 60.0 % of mass 198	32.7	
68	Less than 2.0 % of mass 69	0.0	(0.0)1
69	Mass 69 relative abundance	41.8	
70	Less than 2.0 % of mass 69	0.0	(0.0)1
127	40.0 - 60.0 % of mass 198	58.4	
197	Less than 1.0 % of mass 198	0.2	
198	Base Peak, 100 % relative abundance	100.0	
199	5.0- 9.0 % of mass 198	6.8	
275	10.0 - 30.0 % of mass 198	26.3	
365	Greater than 1.0 % of mass 198	2.9	
441	Present but less than mass 443	11.6	(77.0)3
442	Greater than 40.0 % of mass 198	76.2	
443	17.0 - 23.0 % of mass 442	15.1	(19.8)2

1-Value is % mass 69 2-Value is % mass 442 3-Value is % mass 443

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
	ICIS 180-93544/1	F12250C1.D	12/25/2013	19:23
	IC 180-93544/2	F12250C4.D	12/25/2013	20:50
	IC 180-93544/3	F12250C5.D	12/25/2013	21:18
	IC 180-93544/4	F12250C6.D	12/25/2013	21:46
	IC 180-93544/5	F12250C7.D	12/25/2013	22:15
	IC 180-93544/6	F12250C8.D	12/25/2013	22:43
	IC 180-93544/7	F12250C9.D	12/25/2013	23:11
	IC 180-93544/10	F1225C10.D	12/25/2013	23:41

### FORM VI GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh		ў 	Job No.:	180-28384	384-1				Analy B	Batch No	No.: 93	93544	
SDG No.:		-						:					
Instrument ID: 722		29	Column:	Rxi	-5SilMS	ID:	0.32 (mm)		Heated	Purge:	(Y/N)	Z	
Calibration Start Date: 12/25/2013	19:23	Ca]	alibration	End	Date:	12/25/2013	23:4	⊢:	Calibration	cion ID	13164	64	į
ANALYTE			RRF		. :	CURVE	COEFFICIENT	ENS	# MIN RRF	*RSD #	-	R^2	MIN R^2
	LVL 1	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	TYPE	B M	M2			*RSD	OR COD	OR COD
2-Methylphenol	1.0452	1.0220	1.0631	1.1428	1.0977	Ave	1.0744		0.7000	3.6	20.0		
Indene	1.1042	2.2248	2.2203	2.2382	2.1715	Ave	2.2129		0.0100	5.3	20.0	·	
2,2'-oxybis[1-chloropropane]	0.8815	0.7561	0.7799	0.8076	0.7755	Ave	0.7732	-	0.0100	7.1	20.0	-	
N-Nitrosopyrrolidine	0.5880	0.7252	0.6994	0.5898	0.5678	Ave	0.5709		0.0100	5.5	20.0		
Acetophenone	+++++	1.7081	1.7674	1.8075	1.7310	Ave	1.7146		0.0100	3.9	20.0		
N-Nitrosodi-n-propylamine	0.7291	0.6664	0.7289	0.7603	0.7309	Ave	0.7069		0.5000	5.4	20.0		
Methylphenol, 3 & 4	1.1865	1.0292	1.0817	1.2110	1.1424	Ave	1.1285		0.009	5.3	20.0	 	1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
Hexachloroethane	1.1634 0.5310 0.5605	0.5303	0.5548	0.5645	0.5609	Ave	0.5515		0.3000	2 .4	20.0		
Nitrobenzene	0.2747	0.2608	0.2615	0.2654	0.2748	Ave	0.2645		0.2000	2.8	20.0		
Isophorone	0.4584	0.4187	0.4671	0.4916	0.4925	Ave	0.4713		0.4000	5.2	20.0	:	
2-Nitrophenol	+++++	0.1198	0.1369	0.1653	0.1771	Lin (	0.1235 0.1868		0.1000		20.0	0.9999	0066.0
2,4-Dimethylphenol	0.2490	0.2551	0.2924	0.2968	0.3062	Ave	0.2879	· _	0.2000	7.9	20.0		
Benzoic acid	+++++	+++++	0.0703	0.1336	0.1479	Lin	1.1268 0.2120		0.0100		20.0	0.9972	0.9900
Bis(2-chloroethoxy)methane	0.3462	0.3225	0.3506	0.3294	0.3417	Ave	0.3309		0.3000	4.2	20.0		
2,4-Dichlorophenol	0.2080	0.2213	0.2400	0.2624	0.2804	Ave	0.2542		0.2000	10.9	20.0		manager and the second
1,2,4-Trichlorobenzene	0.3659	0.3309	0.3229	0.3201	0.3329	Ave	0.3287		0.0100	5.0	20.0		
Naphthalene	1.2131	1.0532	1.0758	1.0412	1.0649	Ave	1.0607		0.7000	6.3	20.0		
4-Chloroaniline	0.4370	0.3949	0.4130	0.4431	0.4572	Ave	0.4320		0.0100	4.6	20.0		1
2,6-Dichlorophenol	0.2664	0.2539	0.2704	0.2716	0.2828	Ave	0.2691		0.0100	3.0	20.0		
Hexachlorobutadiene	0.2128	0.1750	0.1747	0.1642	0.1719 Ave	Ave	0.1741		0.0100	9.4	20.0		

Note: The ml coefficient is the same as Ave RRF for an Ave curve type.

# FORM VI GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh		ř 	Job No.:	180-28384-	384-1				Analy		Batch No.:	: 93544	44	
SDG No.:														
Instrument ID: 722		29	Column:		Rxi-5SilMS	ID: 0	.32 (mm)	(2)	Неа	Heated Pu	Purge: (	(X/N)	Z	
Calibration Start Date: 12/25/2013	19:23	Ca]	alibration	End	Date:	12/25/2013		23:41	Cal	Calibration	on ID:	13164	4	a man.
ANALYTE			RRF			CURVE	COE	COEFFICIENT	MIM #	RRF	%RSD #	<u> </u>	-	
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	TYPE		M1 M2				*RSD	OR COD	OR COD
Caprolactam	++++++	+++++	0.0854	0.1016	0.1086	Ave		0.1056	0	0.0100	10.2	20.0		
4-Chloro-3-methylphenol	0.1123	0.1133	0.1124	0.2842	0.2960	Ave		0.2762	0	0.2000 1	11.1	20.0	_	
2-Methylnaphthalene	0.7970	0.7400	0.7538	0.7636	0.7841	Ave		0.7647	-	0.4000	2.5	20.0	:	
1-Methylnaphthalene	0.7959	0.6914	0.7270	0.7106	0.7371	Ave		0.7213	0	0.0100	4.8	20.0		
Hexachlorocyclopentadiene	+ 400 + 600 +	0.2032	0.2440	0.2715	0.2933	Ona 0	0.1233	3.2272 -0.052		0.0500		20.0	0.9999	0.9900
1,2,4,5-Tetrachlorobenzene	0.5990	0.5301	0.5227	0.5051	0.5095	Ave		0.5190	-	0.0100	8.9	20.0	•	
2,4,6-Trichlorophenol	0.3219	0.2380	0.2742	0.3055	0.3226	Ave		0.2919	0	0.2000	14.7	20.0		
2,4,5-Trichlorophenol	++++	0.2736	0.3047	0.3288	0.3364	Ave		0.3277	0	0.2000	8.8	20.0		
1,1'-Biphenyl	1.6600	1.5363	1.5243	1.4592	1,5068	Ave		1.4985	0	0.0100	5.2	20.0		
2-Chloronaphthalene	1.3846	1.2340	1.2211	1.1463	1.1958	Ave		1.1989		0.8000	7.6	20.0		
2-Nitroaniline	1.2034 0.1962 0.2509	1.1264 0.1830 0.2486	0.2210	0.2470	0.2515	Ave		0.2308	0	0.0100	11.9	20.0		
Dimethyl phthalate	1.3310	1.2270	1.2640	1.2418	1.2644 Ave	Ave		1.2587	0	0.0100	2.6	20.0		
1,3-Dinitrobenzene	0.0824	0.1276	0.1589	0.1929	0.2053 Qua		0.1146	4.5643 -0.168		0.0100	<u>:</u>	20.0	0.9998	0.9900
2,6-Dinitrotoluene	0.2111	0.2373	0.2731	0.2921	0.3059	Ave		0.2810		0.2000	13.5	20.0		
Acenaphthylene	1.8344	1.6651	1.8646	1.8634	1.8998	Ave		1.8389		0006.0	4.1	20.0		
3-Nitroaniline	0.2477	0.2487	0.2843	0.3305	0.3444	Ave		0.3143	0	0.0100	14.9	20.0		-
Acenaphthene	1.3498	1.2123	1.2133	1.1896	1.2003	Ave		1.2041	0	0006.0	5.4	20.0		:
2,4-Dinitrophenol	1686	++++	0.0567	0.1142	0.1367	Liņ 1	1.3182 (	0.2092		0.0100		20.0	0.9959	0.9900
4-Nitrophenol	++++	++++	0.1203	0.1395	0.1563	Ave		0.1481	0	0.0100	10.4	20.0		
2,4-Dinitrotoluene	0.2019	0.2767	0.3390	0.3946	0.4099	Lin 0	0.0628	0.4210	0	0.2000		20.0	0.9999	0.9900

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

### FORM VI GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh		dot	b No.:	180-283	8384-1	:				Analy Ba	Batch No	.: 93	544	
SDG No.:											i			
Instrument ID: 722		OS	Column:		Rxi-5SilMS	ID:	0.32 (mm)	u)		Heated F	Purge:	(X/N)	z	
Calibration Start Date: 12/25/2013	19:23	Cal	libration		End Date:	12/25/201	2013	23:41		Calibration	ion ID	: 131	64	
ANALYTE			RRF			CURVE	COE	COEFFICIENT	#	MIN RRF	%RSD #	MAX	R^2 #	MIM
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	TYPE	മ	M1	M2			* KSD	OK COD	OK COD
Dibenzofuran	. (	. 57	- 15	1.6366	1.6802	Ave		1.6808		0.8000	6.3	20.0	-	
-	1.6462	1.6123	1.5893							•	)	) ) 1		
2,3,5,6-Tetrachlorophenol	+++++	0.1907	0.2225	0.2684	0.2886	Qua (	0.1288	3.2891	-0.055	0.0100		20.0	1.0000	0066.0
2,3,4,6-Tetrachlorophenol	+ 5	0.2271	0.2676	0.2946	0.2981	Ave		0.2907		0.0100	11.4	20.0	· · ·	
2-Naphthylamine	+++++	0.5312	0.5619	0.6180	0.6490	Ave		0.5712		0.0100	8.5	20.0		
IN COMMAND AND ADDRESS AND ADD	0.5819	0.5387	0.5174								- 1			
Diethyl phthalate	1.3273	1.2642	1.3560	1.3423	1.3649	Ave		1.3206		0.0100	2.7	20.0	_	
Hexadecane	0.2966	0.3357	0.3919	0.3923	0.3981	Ave		0.3569			10.4			
A O C C C C C C C C C C C C C C C C C C	0.3688	0.3489	0.3231					0.00		000	C	0		
4~Chlorophenyl phenyl ether	0.6447	0.6527	0.6410	0.0401	0.6343	Ave		£/69.0		0.4000	7.0	0.0		
4-Nitroaniline	0.2359	0.2874	0.3342	0.3626	0.3737	Ave		0.3364		0.0100	14.8	20.0		
Fluorene	1.5554	1.3755	1.4200	1.3702	1.4069	Ave		1.4010		0.9000	4.8	20.0		
	1.3812	1.3592	1.3395							i	-			
4,6-Dinitro-2-methylphenol	+ + + + + + + + + + + + + + + + + + + +	0.0431	0.0645	0.1053	0.1167	Lin	0.6766	0.1435		0.0100		20.0	0.9985	0.9900
N-Nitrosodiphenylamine	0.5373	0.4963	0.5186	0.5396	0.5359	Ave		0.5280		0.0100	2.7	20.0		
1,2-Diphenylhydrazine(as Azobenzene)	0.5973	0.5613	0.5990	0.6122	0.5893	Ave		0.5854		0.0100	2.9	20.0		
4-Bromophenyl phenyl ether	0.5778	0.5754	0.5706	0.1948	0.1983	Ave		0.1968		0.1000	4.8	20.0		
	0.1981	0.2041	0.2103								- 1			
Hexachlorobenzene	0.2426	0.2074	0.2111	0.2066	0.2156	Ave		0.2193		0.1000	8.	20.0		
Atrazine	0.1714	0.1216	0.1337	0.1491	0.1519	Ave	<del></del>	0.1429		0.0100	10.7	20.0		
	0.1464	0.1361	0.1331		( ( (		6	0		i c				
Pentachlorophenol	0.0892	0.0972	0.0781	0.1166	0.1209	) Ona	0.2291	7.7808	-0.426	0.0500		70.02	9666.0	0088.0
n-Octadecane	+++++	1.4494	1.7451	1.9845	1.9204	Ave		1.8007			10.2			
Phenanthrene	1.3255	1.1184	1.1215	1.1017	1.1163	Ave		1.1413		0.7000	9.9	20.0	:	
To the transfer of the transfe	1.1171	1.1052	1.1244			f				0	-			
Anthracene	1.1293	1.1258	1.0910	1.1094	1.1368	Ave		1.1116	-	0.007	n.	0.02		
Carbazole	0.9703	0.9224	1.0151	1.0554	1.0725	Ave		1.0318		0.0100	5.7	20.0		

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

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### GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION FORM VI

Lab Name: TestAmerica Pittsburgh		Job	oN d	180-28384	84-1					Analy E	Batch N	oN	3544	
SDG No.:									:					
Instrument ID: 722		SC	Column:	Rxi	-5SilMS	ID: 0	.32 (mm)	(u		Heated	Purge:	(X/N)	z	
Calibration Start Date: 12/25/2013	19:23	Ca	libration	End	Date:	12/25/2013	2013	23:41		Calibration		ID: 13	13164	
ANALYTE			RRF		-	CURVE	COE	COEFFICIENT	=#=	MIN RRF	&RSD	# MAX	R^2 #	
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	되 건 건 건 간		M	M2			₩ K K V	OK COD	 000 000 000 000 000 000 000 000
Di-n-butyl phthalate	0.9509	0.9072	1.0721	1.2532	1.2716	Ave		1.1713		0.0100	14.4	20.0		· i
Fluoranthene	1.0646	0.9807	1.0580	1.1524	1.1872	Ave		1.1369		0.009	8.1	20.0		
Benzidine	+++++	+++++	0.1016	0.1867	0.1977	Lin 0	0.3805	0.2521		0.0100		20.0	0.9993	0.9900
Pyrene	1.1439	1.0315	1.1050	1.1486	1.1962	Ave	:	1.1464		0.6000	4.8	20.0		:
Butyl benzyl phthalate	0.3679	0.3404	0.4031	0.5049	0.5313	Lin 0	0.1190	0.5710		0.0100		20.0	8666.0	0066.0
3,3'-Dichlorobenzidine	0.1711	0.1931	0.2356	0.3237	0.3457	Lin 0	0.2016	0.3975		0.0100		20.0	0.9986	0.9900
Benzo[a]anthracene	1.0220	0.9359	1.0998	1.0699	1.1109	Ave		1.0549	:	0.8000	6.3	20.0		
Bis(2-ethylhexyl) phthalate	0.4156	0.4167	0.5453	0.6865	0.7182	Lin 0	0.0911	0.7512		0.0100		20.0	6666.0	0.9900
Chrysene	1.1007	0.9967	1.0314	0.9839		Ave		1.0115		0.7000	6.6	20.0		
Di-n-octyl phthalate	1.1844	1.5043	0.8476	1.2141	1.3246	Qua 0	0.1340	0.7046 -(	-0.004	0.0100		20.0	19666.0	0.9900
Benzo[b]fluoranthene	1.1887	1.1440	1.1984	1.2652	1,3103	Ave		1.2632		0.7000	0.9	20.0		
7,12-Dimethylbenz(a)anthracene	+++++	0.4524	0.5036	0.5667	0.6153	Ave		0.5843		0.0100	13.7	20.0		
Benzo[k]fluoranthene	1.1925	1.1158	1.1682	1.2119	1.2830	Ave		1.2219		0.7000	4.9	20.0		
Benzo[a]pyrene	0.8966	0.9158	0.9945	1.0670	1.1694	Ave		1.0864		0.7000	12.6	20.0		
Indeno[1,2,3-cd]pyrene	1.1406	1.1128	1.1715	1.2939	1.3727	Ave		1.3027		0.5000	11.1	20.0		
Dibenz(a,h)anthracene	0.9296	0.9619	1.0043	1.0817	1.1678	Ave		1.0941		0.4000	10.6	20.0	:	
Benzo[g,h,i]perylene	1.0422	0.9751	1.0223	1.0815	1.1682	Ave		1.1113		0.5000	8.3	20.0		
2-Fluorophenol (Surr)	0.8920	0.8906	0.8948	0.9973	0.9742	Ave		0.9618			6.1	20.0		
Phenol-d5 (Surr)	1.3327	1.1668	1.3319	1.3219	1.3181	Ave		1.2980			5.3	20.0		
Nitrobenzene-d5 (Surr)	0.2581	0.2381	0.2613	0.2756	0.2791	Ave		0.2665			5.2	20.0		

Note: The ml coefficient is the same as Ave RRF for an Ave curve type.

## Page 1308 of 1943

## GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION FORM VI

Analy Batch No.: 93544	Heated Purge: (Y/N)	Calibration ID: 13164	# MIN RRF 8RSD # MAX R^2 # MIN R^2
	ID: 0.32 (mm)	12/25/2013 23:41	CURVE
Job No.: 180-28384-1	GC Column: Rxi-5SilMS ID: 0.32(mm)	Calibration End Date: 12/25/2013 23:41	RRF
Lab Name: TestAmerica Pittsburgh	Instrument ID: 722	Calibration Start Date: 12/25/2013 19:23	ANALYTE

# MIN R^2			0.9900	
R^2			20.0 0.9998	
MAX Gode	7640	20.0	20.0	20.0
%RSD #		6.1		5.0
MIN RRF %RSD #			0.0100	
#	M2		-1.427	İ
COEFFICIENT	M1	1,3001	0.0989 11.634 -1.427	0.8373
ö	В		0.0989	
CURVE		Ave	Qua	Ave
	LVL 5	1.2808	0.0825 Qua	0.8673
	LVL 4 LVL 5	1.3367 1.2675 1.2808 Ave	0.0784	0.8322 0.8673 Ave
RRF	LVL 3			0.7927
	LVL 1 LVL 2 LVL 6 LVL 7	1.2649 1.2497	0.0380 0.0508	0.8660 0.8701
	LVL 1 LVL 6	1.2649	0.0380	0.8401
ANALYTE		2-Fluorobiphenyl	2,4,6-Tribromophenol (Surr)	Terphenyl-d14 (Surr)

### FORM V GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1 SDG No.:

Lab File ID: F0108DF1.D DFTPP Injection Date: 01/08/2014

DFTPP Injection Time: 01:09 Instrument ID: 722

Analysis Batch No.: 93945

M/E	ION ABUNDANCE CRITERIA		ATIVE DANCE
51	30.0 - 60.0 % of mass 198	38.7	
68	Less than 2.0 % of mass 69	0.0	(0.0)1
69	Mass 69 relative abundance	41.7	
70	Less than 2.0 % of mass 69	0.0	(0.0)1
127	40.0 - 60.0 % of mass 198	55.3	
197	Less than 1.0 % of mass 198	0.0	
198	Base Peak, 100 % relative abundance	100.0	
199	5.0- 9.0 % of mass 198	6.5	-
275	10.0 - 30.0 % of mass 198	25.8	
365	Greater than 1.0 % of mass 198	3.1	
441	Present but less than mass 443	12.9	(79.2)3
442	Greater than 40.0 % of mass 198	88.2	
443	17.0 - 23.0 % of mass 442	16.2	(18.4)2

1-Value is % mass 69 2-Value is % mass 442 3-Value is % mass 443

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 180-93945/2	F01080C1.D	01/08/2014	01:27
SD-191-0-1 DL	180-28384-29 DL	F0108001.D	01/08/2014	02:23

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93945/2 Calibration Date: 01/08/2014 01:27

Instrument ID: 722 Calib Start Date: 12/25/2013 19:23

GC Column: Rxi-5SilMS ID: 0.32(mm) Calib End Date: 12/25/2013 23:41

Lab File ID: F01080C1.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Naphthalene	Ave	1.061	1.063	0.7000	5.01	5.00	0.3	20.0
2-Methylnaphthalene	Ave	0.7647	0.7471	0.4000	4.89	5.00	-2.3	20.0
1-Methylnaphthalene	Ave	0.7213	0.7090	0.0100	4.91	5.00	-1.7	20.0
Acenaphthylene	Ave	1.839	1.934	0.9000	5.26	5.00	5.2	20.0
Acenaphthene	Ave	1.204	1.243	0.9000	5.16	5.00	3.3	20.0
Dibenzofuran	Ave	1.681	1.703	0.8000	5.07	5.00	1.3	20.0
Fluorene	Ave	1.401	1.398	0.9000	4.99	5.00	-0.2	20.0
Phenanthrene	Ave	1.141	1.165	0.7000	5.11	5.00	2.1	20.0
Anthracene	Ave	1.112	1.193	0.7000	5.37	5.00	7.3	20.0
Carbazole	Ave	1.032	1.102	0.0100	5.34	5.00	6.8	20.0
Fluoranthene	Ave	1.137	1.175	0.6000	5.17	5.00	3.3	20.0
Pyrene	Ave	1.146	1.125	0.6000	4.91	5.00	-1.9	20.0
Benzo[a]anthracene	Ave	1.055	1.047	0.8000	4.96	5.00	-0.8	20.0
Chrysene	Ave	1.012	0.9857	0.7000	4.87	5.00	-2.6	20.0
Benzo[b]fluoranthene	Ave	1.263	1.226	0.7000	4.85	5.00	-3.0	20.0
Benzo[k]fluoranthene	Ave	1.222	1.261	0.7000	5.16	5.00	3.2	20.0
Benzo[a]pyrene	Ave	1.086	1.080	0.7000	4.97	5.00	-0.6	20.0
Indeno[1,2,3-cd]pyrene	Ave	1.303	1.309	0.5000	5.02	5.00	0.5	20.0
Dibenz(a,h)anthracene	Ave	1.094	1.114	0.4000	5.09	5.00	1.8	20.0
Benzo[g,h,i]perylene	Ave	1.111	1.120	0.5000	5.04	5.00	0.8	20.0
2-Fluorophenol (Surr)	Ave	0.9618	1.159		6.03	5.00	20.5*	20.0
Phenol-d5 (Surr)	Ave	1.298	1.527		5,88	5.00	17.6	20.0
Nitrobenzene-d5 (Surr)	Ave	0.2665	0.2998		5.62	5.00	12.5	20.0
2-Fluorobiphenyl	Ave	1.300	1.326		5.10	5.00	2.0	20.0
2,4,6-Tribromophenol (Surr)	Qua	0.0745	0.0831	0.0100	5.17	5.00	3.4	20.0
Terphenyl-d14 (Surr)	Ave	0.8373	0.8262		4.93	5.00	-1.3	20.0

#### FORM V

### GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

DFTPP Injection Date: 12/11/2013 Lab File ID: D1211DF1.D

Instrument ID: 732 DFTPP Injection Time: 05:18

Analysis Batch No.: 92134

M/E	ION ABUNDANCE CRITERIA		LATIVE DANCE
51	30.0 - 60.0 % of mass 198	38.5	
68	Less than 2.0 % of mass 69	0.0	(0.0)1
69	Mass 69 relative abundance	41.0	
70	Less than 2.0 % of mass 69	0.0	(0.0)1
127	40.0 - 60.0 % of mass 198	42.6	-
197	Less than 1.0 % of mass 198	0.0	
198	Base Peak, 100 % relative abundance	100.0	
199	5.0- 9.0 % of mass 198	7.0	
275	10.0 - 30.0 % of mass 198	27.0	
365	Greater than 1.0 % of mass 198	2.3	
441	Present but less than mass 443	2.9	(25.9)3
442	Greater than 40.0 % of mass 198	58.1	
443	17.0 - 23.0 % of mass 442	11.4	(19.6)2

1-Value is % mass 69 2-Value is % mass 442 3-Value is % mass 443

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 180-92134/12	D1211IC1.D	12/11/2013	05:33
	IC 180-92134/2	D1211IC2.D	12/11/2013	05:59
	IC 180-92134/3	D1211IC3.D	12/11/2013	06:26
	ICIS 180-92134/4	D1211IC4.D	12/11/2013	06:52
	IC 180-92134/5	D1211IC5.D	12/11/2013	07:18
	IC 180-92134/6	D1211IC6.D	12/11/2013	07:45
	IC 180-92134/7	D1211IC7.D	12/11/2013	08:11
	IC 180-92134/8	D1211IC8.D	12/11/2013	08:37

## GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION FORM VI

Lab Name: TestAmerica Pittsburgh		dob	oN o	180-28384-1	84-1					Analy B	Batch No.		92134	
Instrument ID: 732		25	Column:	Rxi	-5SilMS	ID:	0.32 (mm)	(1		Heated	Purge:	(X/N)	Z	:
Calibration Start Date: 12/11/2013	05:33	Cal	ibration	End	Date:	12/11/2013	İ	08:37		Calibration		ID: 12823	323	
ANALYTE			RRE			CURVE	COE	COEFFICIENT		# MIN RRF	%RSD	# MAX	R^2	# MIN R^2
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	E A A A	т	M	M2			** XX VV	OK COD	OK COL
2-Methylphenol	0.9219	1.0613	1.1599	1.1805	1.2010	Ave		1.1324		0.7000	8.4	20.0		
Indene	2.0902	2.2738	2.2807	2.2763	2.3216	Ave		2.2535		0.0100	3.1	20.0		
2,2'-oxybis[1-chloropropane]	1.5649	1.6438	1.6760	1.6245	1.6020	Ave		1.5839		0.0100	4.2	20.0	:	
N-Nitrosopyrrolidine	+ 00 +	0.4404	0.5038	0.5412	0.5778	Ave		0.5434		0.0100	10.0	20.0		:
Acetophenone	1.7204	1.7069	1.8264	1.7662	1.7963	Ave		1.7487		0.0100	2.9	20.0		
N-Nitrosodi-n-propylamine	0.6984	0.8663	0.8859	0.9072	0.8997	Ave		0.8499		0.5000	7.8	20.0		
Methylphenol, 3 & 4	1.0361	1.1798	1.2616	1.2460	1.2920	Ave		1.2230		0.009	9.9	20.0		
Hexachloroethane		0.5617	0.5643	0.5682	0.5823	Ave		0.5685		0.3000	1.4	20.0		
Nitrobenzene	1	0.3165	0.3260	0.3198	0.3281	Ave		0.3169		0.2000	7.3	20.0		
Isophorone	0.4068	0.5255	0.5505	0.5615	0.5824	Ave		0.5497		0.4000	11.8	20.0		
2-Nitrophenol	0.5901	0.5910 0.1451 0.1992	0.6037 0.1670 0.2028	0.1777	0.1883	Ave		0.1821		0.1000	11.3	20.0		i
2,4-Dimethylphenol	0.2407	0.2930	0.3235	0.3158	0.3241	Ave		0.3096		0.2000	9.7	20.0	:	-
Benzoic acid	0.2074	0.1226	0.1244	0.1564	0.1818	Lin (	0.4901 (	0.2229		0.0100		20.0	0.9977	0.9900
Bis(2-chloroethoxy)methane	0.3331	0.3691	0.3645	0.3603	0.3604	Ave		0.3598		0.3000	3.1	20.0		<u> </u>
2,4-Dichlorophenol	0.2347	0.2862	0.3045	0.3034	0.3147	Ave		0.3003		0.2000	9.7	20.0		
1,2,4-Trichlorobenzene	0.3479	0.3587	0.3710	0.3559	0.3598	Ave		0.3574		0.0100	2.1	20.0		
Naphthalene	1.0584	1.0406	1.0573	1.0323	1.0360	Ave	····	1.0482		0.7000	1.5	20.0	İ	
4-Chloroaniline	0.3799	0.4021	0.4214	0.4295	0.4323	Ave	. — . — .	0.4229		0.0100	5.	20.0		
2,6-Dichlorophenol	0.2552	0.2907	0.2963	0.3036	0908.0	Ave		0.2976		0.0100	6.2	20.0		!
Hexachlorobutadiene	0.2202	0.2220	0.2299	0.2173	0.2209	Ave		0.2220		0.0100	1.7	20.0		

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

## FORM VI GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh		dot	op No.:	180-28384-	384-1				Analy E	Analy Batch No.		92134	
SDG No.:	:									:		:	
Instrument ID: 732		29	Column:	Rxi	-5SilMS	ID:	0.32 (mm)	(1	Heated	Purge:	(Y/N)	Z	
Calibration Start Date: 12/11/2013	05:33	Cal	libration	on End	Date:	12/11/2013		08:37	Calibration		ID: 12823	323	
ANALYTE			RRF			CURVE	COE	COEFFICIENT	# MIN RRF	%RSD	# MAX	R^2 #	MIN R^2
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	I	т	M1 M2	-		*RSD	OK COD	OK COD
Caprolactam	0.0413	0.0624	0.0737	0.0857	0.0943	Lin	0.1694 (	0.1057	0.0100		20.0	0.9990	0.9900
4-Chloro-3-methylphenol	0.1005	0.1012	0.1058	0.2896	0.2983	Ave		0.2842	0.2000	9.3	20.0		
2-Methylnaphthalene	0.6962	0.7351	0.3079	0.7357	0.7410 Ave	Ave		0.7405	0.4000	2.7	20.0		
1-Methylnaphthalene	0.6390	0.6819	0.7061	0.6835	0.6867	Ave		0.6875	0.0100	3.2	20.0		
Hexachlorocyclopentadiene	0.3004	0.3318	0.3528	0.3868	0.4254	Ave		0.3846	0.0500	13.1	20.0		
1,2,4,5-Tetrachlorobenzene	0.6497	0.6491	0.6414	0.6459	0.6524	Ave		0.6357	0.0100	2.9	20.0		
2,4,6-Trichlorophenol	0.2436	0.3265	0.3508	0.3758	0.3842	Ave	-	0.3537	0.2000	13.9	20.0		:
2,4,5-Trichlorophenol	0.2767	0.3465	0.3963	0.3885	0.4156	Ave		0.3801	0.2000	12.4	20.0		
1,1'-Biphenyl	1.4314	1.4528	1.4915	1.4358	1.5040	Ave		1.4485	0.0100	2.2	20.0	:	
2-Chloronaphthalene	1.0882	1.2539	1.2284	1.1933	1.1641	Ave		1.1754	0.8000	4.5	20.0		
2-Nitroaniline	+++++	0.2111	0.2484	0.2854	0.3045	Ave		0.2810	0.0100	13.3	20.0		
Dimethyl phthalate	1.2811	1.2509	1.2954	1.2388	1.3192	Ave		1.2585	0.0100	3.9	20.0		-
1,3-Dinitrobenzene	+++++	0.1341	0.1666	0.1911	0.2158	Lin	0.1481 (	0.2247	0.0100		20.0	0.9997	0.9900
2,6-Dinitrotoluene	+ + + + + + + + + + + + + + + + + + + +	0.2568	0.2631	0.2830	0.2933	Ave		0.2832	0.2000	5.9	20.0		
Acenaphthylene	1.8301	1.7266	1.8341	1.7726	1.8648	Ave		1.7599	0.9000	6.9	20.0		
3-Nitroaniline	+++++	0.2548	0.2770	0.3080	0.3359	Ave	:	0.3089	0.0100	0 10.2	20.0		
2,4-Dinitrophenol	+++++	0.1099	0.1127	0.1586	0.2016	Lin	0.7013	0.2463	0.0100		20.0	0.9982	0.9900
Acenaphthene	1,1791	1,1816	1,2161	1.1899	1,2041	Ave		1.1759	0.000.0	2.4	20.0		
4-Nitrophenol	+++++	0.1282	0.1347	0.1487	0.1578	Ave		0.1479	0.0100	0.7	20.0		
2,4-Dinitrotoluene	<u>.:</u>	0.3336	0.3584	0.3832	0.4193	Ave		0.3894	0.2000	8.4	20.0		

Note: The ml coefficient is the same as Ave RRF for an Ave curve type.

### GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION FORM VI

Lab Name: TestAmerica Pittsburgh		Job	ob No.:	180-28384	84-1			Analy B	Batch No	No.: 92	92134	
SDG No.:				i					:			
Instrument ID: 732		SS	Column:	Rxi	-5SilMS ID:	0.32 (mm)		Heated	Purge:	(X/N)	Z	
Calibration Start Date: 12/11/2013	05:33	Cal	alibration	End	Date: 12/1	12/11/2013 0	08:37	Calibration	tion ID	): 1282	23	
ANALYTE		:	RRF		CURVE	:	COEFFICIENT	# MIN RRE	%RSD #		R^2	MIN R^2
	LVL 1	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5 TYPE	<b>.</b>	M1 M2	· n· · = ···		%RSD	OR COD	OR COD
Dibenzofuran	(1) (r	1.7343	1.7203	1.6779	1.7168 Ave		1.6769	0.8000	3.3	20.0		
2,3,5,6-Tetrachlorophenol	++++	0.3061	0.3438	0.3683	0.3934 Ave	0	0.3708	0.0100	9.3	20.0		
2,3,4,6-Tetrachlorophenol	0.2714	0.3467	0.3715	0.3800	0.3981 Ave	0	0.3671	0.0100	11.4	20.0		
2-Naphthylamine	0.4470	0.5771	0.5698	0.5937	0.6111 Ave	0	0.5439	0.0100	10.3	20.0		
Diethyl phthalate	1.1361	1.1933	1.2252	1.2265	1.2599 Ave		1.2086	0.0100	3.1	20.0		
Hexadecane	0.3704	0.4425	0.4583	0.4574	0.4565 Ave	0	0.4420		6.7			
4-Chlorophenyl phenyl ether	0.7333	0.6996		0.7215	0.7474 Ave	0	0.7231	0.4000	2.0	20.0		
4-Nitroaniline	+++++	0.2604	0.2932	0.3194	0.3459 Ave	0	0.3179	0.0100	9.7	20.0		
Fluorene	1,2096	1.3765	1.3887	1.3618	1.4096 Ave		1.3507	0.9000	4.5	20.0		
4,6-Dinitro-2-methylphenol	+++++	0.0833	0.1124	0.1339	0.1487 Lin	0.4597 0	0.1693	0.0100		20.0	0.9994	0.9900
N-Nitrosodiphenylamine	0.4707	0.5226	0.5287	0.5348	0.5423 Ave	0	0.5278	0.0100	4.6	20.0		
1,2-Diphenylhydrazine(as Azobenzene)	0.5120	0.6144	0.6235	0.6379	0.6336 Ave	0	0.6173	0.0100	7.1	20.0		
4-Bromophenyl phenyl ether	0.2048	0.2219	0.2186	0.2245	0.2331 Ave	0	0.2260	0.1000	4.8	20.0		
Hexachlorobenzene	0.2311	0.2221	0.2168	0.2125	0.2205 Ave	0	0.2200	0.1000	2.4	20.0		
Atrazine	++++	0.1468	0.1574	0.1684	0.1709 Qua	0.0703 5	5.3037 0.8983	0.0100		20.0	6666.0	0.066.0
Pentachlorophenol	0.2837	0.1695	0.1167	0.1374	0.1530 Lin	0.2349 0	0.1671	0.0500		20.0	0.9991	0.9900
n-Octadecane	1.3147	1.7186	1.9051	1.9668	1.9485 Ave		1.8196		12.0	-		
Phenanthrene	1.1071	1.1188	1.1172	1.1019	1.1074 Ave		1.1101	0.7000	0	20.0	-	
Anthracene	0.9659	1.1137	1.1063	1.1328	1.1480 Ave		1.1129	0.7000	5.5	20.0		
Carbazole	0.8440	0.9812	0.9863	1.0406	1.0461 Ave	Г	1.0056	0.0100	7.1	20.0		

Note: The ml coefficient is the same as Ave RRF for an Ave curve type.

## FORM VI GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh		Job	b No.:	180-28384-	384-1					Analy Batch No.:	ch No.	: 92134	4	
SDG No.:					:									
Instrument ID: 732		SB	Column:	Rx1-	5SilMS	: ID:	0.32 (mm)		!	Heated P	Purge: (	(Y/N) N		
Calibration Start Date: 12/11/2013	05:33	Cal	libration	End	Date:	12/11/201	m	08:37		Calibration	on ID:	12823		,
ANALYTE			RRF			CURVE	COE	COEFFICIENT	***	MIN RRE	%RSD #		R^2 #	MIN R^2
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	TYPE	щ	M1	M2			%RSD OF	OR COD	OR COD
Di-n-butyl phthalate	+ 0	0.9127	0.9921	1.0963	1.1513	Ave		1.1056		0.0100	10.3	20.0		
Fluoranthene	1.0377	1.2024	1.2199	1.2806	1.3359	Ave		1.2582		0.6000	8.2	20.0		
Benzidine	+++++	++++	0.1349	0.1770	0.2799	Qua	0.3228	3,3992	-0.050	0.0100		20.00	0.9987	0066.0
Pyrene	0.9336	1.0394	1.0823	1.0965	1.1325	Ave		1.0812		00009.0	6.2	20.0		
Butyl benzyl phthalate	0.2120	0.2835	0.3308	0.3851	0.4172	Lin	0.1215 (	0.4457	-	0.0100		20.0	0.9995	0066.0
3,3'-Dichlorobenzidine	0.1618	0.2375	0.2628	0.3237	0.3823	Qua	0.1204	2.5434	-0.039	0.0100		20.0	8666.0	0066.0
Bis(2-ethylhexyl) phthalate	0.2497	0.3676	0.4149	0.5202	0.5738	Lin	0.1365 (	0.6204		0.0100	i	20.00	7666.0	0.9900
Benzo[a]anthracene	0.5969	1.0416	1.0850	1.0942	1.1048	Ave		1.0792		0.8000	3.7	20.0		
Chrysene	0.9588	1.0155	1.0149	1.0284	1.0385	Ave		1.0209		0.7000	2.7	20.0		
Di-n-octyl phthalate	1,1618	0.5739	0.6792	0.8661	1.0536	Qua	0.2335 (	0.8397	-0.002	0.0100		20.0	9666.0	0.9900
7,12-Dimethylbenz(a)anthracene	0.4171	0.4839	0.5492	0.5849	0.6299	Ave		0.5682		0.0100	14.1	20.0	-	·-·
Benzo[b]fluoranthene	0.9866	1.1502	1.2423	1.2359	1.3244	Ave		1.2319		0.7000	9.3	20.0		
Benzo[k]fluoranthene	1.1761	1.2409	1.2555	1.3540	1.3118	Ave		1.2783		0.7000	4.3	20.0		
Benzo[a]pyrene	0.8615	0.9745	1.0436	1.1059	1.1896	Ave		1.0950		0.7000	11.5	20.0		
Indeno[1,2,3-cd]pyrene	0.8670	1.0470	1.1175	1.1871	1.2414	Ave		1.1602	-	0.5000	12.5	20.0		:
Dibenz(a,h)anthracene	0.7091	0.8827	0.9628	1.0063	1.0323	Ave		0.9749		0.4000	12.9	20.0		
Benzo[g,h,i]perylene	0.7731	0.9034	0.9915	1.0207	1.0456	Ave		0.9956		0.5000	10.8	20.0		
2-Fluorophenol (Surr)	0.8733	1.0165	1.0587	1.0642	1.1345	Ave		1.0741			6.8	20.0		
Phenol-d5 (Surr)	1.1126	1.3138	1.4230	1.4501	1.5074	Ave		1.4153			6.6	20.0		
Nitrobenzene-d5 (Surr)	0.2354	0.2855	0.3052	0.3134	0.3179	Ave		0.3040	-		10.1	20.0		

Note: The ml coefficient is the same as Ave RRF for an Ave curve type.

## GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION FORM VI

	:		
Analy Batch No.: 92134		Heated Purge: (Y/N) N	Calibration ID: 12823
Job No.: 180-28384-1		GC Column: Rxi-5SilMS ID: 0.32(mm)	Calibration End Date: 12/11/2013 08:37
Lab Name: TestAmerica Pittsburgh	SDG No.:	Instrument ID: 732	Calibration Start Date: 12/11/2013 05:33

ANALYTE			RRE			CURVE	20	COEFFICIENT	-	MIM #	MIN RRF %	%RSD #	MAX	R^2	# MIN R^2	MIN R^2
	LVL 1 LVL 2 LVL 6 LVL 7	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	권 건 건	ш	M1	M2				9.R30	OK (	5	
2-Fluorobiphenyl	1.3056 1.3765	1.3765	1.3781	1.3468	1.3946 Ave	Ave		1.3463				2.5	20.0			
2,4,6-Tribromophenol (Surr)	1.3360 1.3168	0.0672	0.0766	0.0816	0.0866 Ave	Ave		0.0838		0.0	0.0100 11.3	1.3	20.0			
	0.0893 0.0917	0.0917	0.0940													
Terphenyl-d14 (Surr)	0.6985	0.7543	0.7950		0.8083 0.8350 Ave	Ave		0.8022			-	6.5	20.0	-		
	0.8427 0.8316	0.8316	0.8525	1				_				_		_		

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#### FORM V

### GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

Lab File ID: D1231DF1.D DFTPP Injection Date: 12/31/2013

Instrument ID: 732 DFTPP Injection Time: 10:16

Analysis Batch No.: 93613

M/E	ION ABUNDANCE CRITERIA		ATIVE DANCE
51	30.0 - 60.0 % of mass 198	40.9	
68	Less than 2.0 % of mass 69	0.0	(0.0)1
69	Mass 69 relative abundance	42.6	
70	Less than 2.0 % of mass 69	0.1	(0.2)1
127	40.0 - 60.0 % of mass 198	43.3	
197	Less than 1.0 % of mass 198	0.0	
198	Base Peak, 100 % relative abundance	100.0	
199	5.0- 9.0 % of mass 198	6.6	
275	10.0 - 30.0 % of mass 198	27.6	
365	Greater than 1.0 % of mass 198	2.6	
441	Present but less than mass 443	6.0	(55.8)3
442	Greater than 40.0 % of mass 198	55.6	
443	17.0 - 23.0 % of mass 442	10.7	(19.3)2

1-Value is % mass 69 2-Value is % mass 442 3-Value is % mass 443

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 180-93613/22	D12310CC.D	12/31/2013	10:31
	MB 180-93179/1-A	D1231003.D	12/31/2013	10:59
	LCS 180-93179/2-A	D1231004.D	12/31/2013	11:51
SD-163-0-1 MS	180-28384-1 MS	D1231005.D	12/31/2013	14:03
SD-163-0-1 MSD	180-28384-1 MSD	D1231006.D	12/31/2013	14:29
SD-163-0-1	180-28384-1	D1231007.D	12/31/2013	14:56
SD-164-0-1	180-28384-2	D1231008.D	12/31/2013	15:22
SD-165-0-1	180-28384-3	D1231009.D	12/31/2013	15:49
SD-166-0-1	180-28384-4	D1231010.D	12/31/2013	16:15
SD-167-0-1	180-28384-5	D1231011.D	12/31/2013	16:41
SD-168-0-1	180-28384-6	D1231012.D	12/31/2013	17:08
SD-169-0-1	180-28384-7	D1231013.D	12/31/2013	17:34
SD-170-0-1	180-28384-8	D1231014.D	12/31/2013	18:00
SD-171-0-1	180-28384-9	D1231015.D	12/31/2013	18:26
SD-172-0-3	180-28384-10	D1231016.D	12/31/2013	18:53
SD-173-0-3	180-28384-11	D1231017.D	12/31/2013	19:19
SD-174-0-1	180-28384-12	D1231018.D	12/31/2013	19:46
SD-175-0-1	180-28384-13	D1231019.D	12/31/2013	20:12
SD-176-0-3	180-28384-14	D1231020.D	12/31/2013	20:38
SD-177-0-2	180-28384-15	D1231021.D	12/31/2013	21:04
SD-178-0-1	180-28384-16	D1231022.D	12/31/2013	21:31
SD-179-0-1	180-28384-17	D1231023.D	12/31/2013	21:57

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93613/22 Calibration Date: 12/31/2013 10:31

Instrument ID: 732 Calib Start Date: 12/11/2013 05:33

GC Column: Rxi-5SilMS ID: 0.32(mm) Calib End Date: 12/11/2013 08:37

Lab File ID: D12310CC.D Conc. Units: ng/uL

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ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,4-Dioxane	Ave	0.4034	0.4052	0.0100	5.02	5.00	0.4	20.0
N-Nitrosodimethylamine	Ave	0.5369	0.5861	0.0100	5.46	5.00	9.2	20.0
Pyridine	Ave	. 0.9657	1.036	0.0100	5,37	5.00	7.3	20.0
Methyl methanesulfonate	Ave	0.5820	0.6183	0.0100	5.31	5.00	6.2	20.0
Benzaldehyde	Qua	0.6652	0.8454	0.0100	4.93	5.00	-1.4	20.0
Aniline	Ave	1.715	1.863	0.0100	5.43	5.00	8.7	20.0
Phenol	Ave	1.632	1.679	0.8000	5.14	5.00	2.9	20.0
Bis(2-chloroethyl)ether	Ave	1.106	1.123	0.7000	5.08	5.00	1.5	20.0
2-Chlorophenol	Ave	1.282	1.349	0.8000	5.26	5.00	5.3	20.0
1,3-Dichlorobenzene	Ave	1.589	1.575	0.0100	4.96	5.00	-0.9	20.0
1,4-Dichlorobenzene	Ave	1.610	1.620	0.0100	5.03	5.00	0.6	20.0
Benzyl alcohol	Ave	0.7606	0.7499	0.0100	4.93	5.00	-1.4	20.0
1,2-Dichlorobenzene	Ave	1.532	1.549	0.0100	5.05	5.00	1.1	20.0
2-Methylphenol	Ave	1.132	1.186	0.7000	5.24	5.00	4.7	20.0
Indene	Ave	2.253	2.305	0.0100	5.11	5.00	2.3	20.0
2,2'-oxybis[1-chloropropane]	Ave	1.584	1.648	0.0100	5.20	5.00	4.0	20.0
N-Nitrosopyrrolidine	Ave	0.5434	0.5660	0.0100	5.21	5.00	4.2	20.0
Acetophenone	Ave	1.749	1.824	0.0100	5.22	5.00	4.3	20.0
N-Nitrosodi-n-propylamine	Ave	0.8499	0.9138	0.5000	5.38	5.00	7.5	20.0
Methylphenol, 3 & 4	Ave	1.223	1.264	0.6000	5.17	5.00	3.4	20.0
Hexachloroethane .	Ave	0.5685	0.5937	0.3000	5.22	5.00	4.4	20.0
Nitrobenzene	Ave	0.3169	0.3312	0.2000	5.23	5.00	4.5	20.0
Isophorone	Ave	0.5497	0.5728	0.4000	5.21	5.00	4.2	20.0
2-Nitrophenol	Ave	0.1821	0.1846	0.1000	5.07	5.00	1.4	20.0
2,4-Dimethylphenol	Ave	0.3096	0.3140	0.2000	5.07	5.00	1.4	20.0
Benzoic acid	Lin	0.1791	0.1532	0.0100	8.83	10.0	-11.7	20.0
Bis(2-chloroethoxy)methane	Ave	0.3598	0.3547	0.3000	4.93	5.00	-1.4	20.0
2,4-Dichlorophenol	Ave	0.3003	0.3053	0.2000	5.08	5.00	1.7	20.0
1,2,4-Trichlorobenzene	Ave	0.3574	0.3514	0.0100	4.92	5.00	-1.7	20.0
Naphthalene	Ave	1.048	1.030	0.7000	4.91	5.00	-1.7	20.0
4-Chloroaniline	Ave	0.4229	0.4102	0.0100	4.85	5.00	-3.0	20.0
2,6-Dichlorophenol	Ave	0.2976	0.3084	0.0100	5.18	5.00	3.6	20.0
Hexachlorobutadiene	Ave	0.2220	0.2207	0.0100	4.97	5.00	-0.6	20.0
Caprolactam	Lin	0.0831	0.0855	0.0100	4.72	5.00	-5.6	20.0
4-Chloro-3-methylphenol	Ave	0.2842	0.2849	0.2000	5.01	5.00	0.3	20.0
2-Methylnaphthalene	Ave	0.7405	0.7220	0.4000	4.87	5.00	-2.5	20.0
1-Methylnaphthalene	Ave	0.6875	0.6750	0.0100	4.91	5.00	-1.8	20.0
Hexachlorocyclopentadiene	Ave	0.3846	0.3780	0.0500	4.91	5.00	-1.7	20.0
1,2,4,5-Tetrachlorobenzene	Ave	0.6357	0.6417	0.0100	5.05	5.00	0.9	20.0
2,4,6-Trichlorophenol	Ave	0.3537	0.3736	0.2000	5.28	5.00	5.6	20.0
2,4,5-Trichlorophenol	Ave	0.3801	0.4034	0.2000	5.31	5.00	6.1	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93613/22 Calibration Date: 12/31/2013 10:31

Instrument ID: 732 Calib Start Date: 12/11/2013 05:33

GC Column: Rxi-5SilMS ID: 0.32(mm) Calib End Date: 12/11/2013 08:37

Lab File ID: D12310CC.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,1'-Biphenyl	Ave	1.449	1.428	0.0100	4.93	5.00	-1.4	20.0
2-Chloronaphthalene	Ave	1.175	1.153	0.8000	4.91	5.00	-1.9	20.0
2-Nitroaniline	Ave	0.2810	0.2997	0.0100	5.33	5.00	6.7	20.0
Dimethyl phthalate	Ave	1.258	1.245	0.0100	4.95	5.00	-1.1	20.0
1,3-Dinitrobenzene	Lin	0.1947	0.2039	0.0100	5.13	5.00	2.6	20.0
2,6-Dinitrotoluene	Ave	0.2832	0.2799	0.2000	4.94	5.00	-1.2	20.0
Acenaphthylene	Ave	1.760	1.801	0.9000	5.12	5.00	2.3	20.0
3-Nitroaniline	Ave	0.3089	0.2996	0.0100	4.85	5.00	-3.0	20.0
2,4-Dinitrophenol	Lin	0.1825	0.1715	0.0100	9.77	10.0	-2.3	20.0
Acenaphthene	Ave	1.176	1.169	0.9000	4.97	5.00	-0.6	20.0
4-Nitrophenol	Ave	0.1479	0.1560	0.0100	10.6	10.0	5.5	20.0
2,4-Dinitrotoluene	Ave	0.3894	0.3797	0.2000	4.88	5.00	-2.5	20.0
Dibenzofuran	Ave	1.677	1.645	. 0.8000	4.91	5.00	-1.9	20.0
2,3,5,6-Tetrachlorophenol	Ave	0.3708	0.3599	0.0100	4.85	5.00	-2.9	20.0
2,3,4,6-Tetrachlorophenol	Ave	0.3671	0.3545	0.0100	4.83	5.00	-3.4	20.0
2-Naphthylamine	Ave	0.5439	0.5875	0.0100	5.40	5.00	8.0	20.0.
Diethyl phthalate	Ave	1.209	1.277	0.0100	5.28	5.00	5.7	20.0
4-Chlorophenyl phenyl ether	Ave	0.7231	0.7146	0.4000	4.94	5.00	-1.2	20.0
4-Nitroaniline	Ave	0.3179	0.3194	0.0100	5.02	5.00	0.5	20.0
Fluorene	Ave	1.351	1.332	0.9000	4.93	5.00	-1.4	20.0
4,6-Dinitro-2-methylphenol	Lin	0.1380	0.1445	0.0100	10.4	10.0	3.7	20.0
N-Nitrosodiphenylamine	Ave	0.5278	0.5406	0.0100	5.12	5.00	2.4	20.0
1,2-Diphenylhydrazine(as Azobenzene)	Ave	0.6173	0.6502	0.0100	5.27	5.00	5.3	20.0
4-Bromophenyl phenyl ether	Ave	0.2260	0.2318	0.1000	5.13	5.00	2.6	20.0
Hexachlorobenzene	Ave	0.2200	0.2170	0.1000	4.93	5.00	-1.4	20.0
Atrazine	Qua	0.1590	0.1703	0.0100	4.96	5.00	-0.8	20.0
Pentachlorophenol	Lin	0.1688	0.1553	0.0500	10.2	10.0	2.3	20.0
Phenanthrene	Ave	1.110	1.097	0.7000	4.94	5.00	-1.1	20.0
Anthracene	Ave	1,113	1.160	0.7000	5.21	5.00	4.3	20.0
Carbazole	Ave	1.006	1.037	0.0100	5.15	5.00	3.1	20.0
Di-n-butyl phthalate	Ave	1.106	1.128	0.0100	5.10	5.00	2.0	20.0
Fluoranthene	Ave	1.258	1.293	0.6000	5.14	5.00	2.8	20.0
Benzidine	Qua	0.2433	0.2119	0.0100	4.88	5.00	-2.4	20.0
Pyrene	Ave	1.081	1.129	0.6000	5.22	5.00	4.4	20.0
Butyl benzyl phthalate	Lin	0.3673	0.4090	0.0100	5.07	5.00	1.5	20.0
3,3'-Dichlorobenzidine	Qua	0.3234	0.3499	0.0100	4.90	5.00	-2.0	20.0
Bis(2-ethylhexyl) phthalate	Lin	0.4963	0.5715	0.0100	5.15	5.00	3.0	20.0
Benzo[a]anthracene	Ave	1.079	1.088	0.8000	5.04	5.00	0.8	20.0
Chrysene	Ave	1.021	0.998	0.7000	4.89	5.00	-2.2	20.0
Di-n-octyl phthalate	Qua	0.9594	1.112	0.0100	5.59	5.00	11.7	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93613/22 Calibration Date: 12/31/2013 10:31

Instrument ID: 732 Calib Start Date: 12/11/2013 05:33

Lab File ID: D12310CC.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
7,12-Dimethylbenz(a)anthrace	Ave	0.5682	0.5778	0.0100	5.08	5.00	1.7	20.0
Benzo[b]fluoranthene	Ave	1.232	1.292	0.7000	5.25	5.00	4.9	20.0
Benzo[k]fluoranthene	Ave	1.278	1.227	0.7000	4.80	5.00	-4.0	20.0
Benzo[a]pyrene	Ave	1.095	1.156	0.7000	5.28	5.00	5.5	20.0
Indeno[1,2,3-cd]pyrene	Ave	1.160	1.229	0.5000	5.30	5.00	6.0	20.0
Dibenz(a,h)anthracene	Ave	0.9749	1.035	0.4000	5.31	5.00	6.2	20.0
Benzo[g,h,i]perylene	Ave	0.996	1.057	0.5000	5.31	5.00	6.1	20.0
2-Fluorophenol (Surr)	Ave	1.074	1.132		5.27	5.00	5.4	20.0
Phenol-d5 (Surr)	Ave	1.415	1.511		5.34	5.00	6.8	20.0
Nitrobenzene-d5 (Surr)	Ave	0.3040	0.3313		5.45	5.00	9.0	20.0
2-Fluorobiphenyl	Ave	1.346	1.339		4.97	5.00	-0.6	20.0
2,4,6-Tribromophenol (Surr)	Ave	0.0838	0.0878	0.0100	5.23	5.00	4.7	20.0
Terphenyl-dl4 (Surr)	Ave	0.8022	0.8376		5.22	5.00	4.4	20.0

### FORM V GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK

DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab File ID: D0102DF1.D DFTPP Injection Date: 01/02/2014

Instrument ID: 732 DFTPP Injection Time: 12:00

Analysis Batch No.: 93710

M/E	ION ABUNDANCE CRITERIA		ATIVE DANCE
51	30.0 - 60.0 % of mass 198	42.0	
68	Less than 2.0 % of mass 69	0.0	(0.0)1
69	Mass 69 relative abundance	43.5	
70	Less than 2.0 % of mass 69	0.0	(0.0)1
127	40.0 - 60.0 % of mass 198	44.8	
197	Less than 1.0 % of mass 198	0.0	-
198	Base Peak, 100 % relative abundance	100.0	
199	5.0- 9.0 % of mass 198	6.9	
275	10.0 - 30.0 % of mass 198	26.4	
365	Greater than 1.0 % of mass 198	2.1	
441	Present but less than mass 443	8.6	(86.3)3
442	Greater than 40.0 % of mass 198	53.5	
443	17.0 - 23.0 % of mass 442	10.0	(18.7)2

1-Value is % mass 69

2-Value is % mass 442

3-Value is % mass 443

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 180-93710/10	D01020CC.D	01/02/2014	12:15
SD-180-0-1	180-28384-18	D0102001.D	01/02/2014	19:05
SD-181-0-1	180-28384-19	D0102002.D	01/02/2014	19:32
SD-182-0-3	180-28384-20	D0102003.D	01/02/2014	19:58

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93710/10 Calibration Date: 01/02/2014 12:15

Instrument ID: 732 Calib Start Date: 12/11/2013 05:33

Lab File ID: D01020CC.D Conc. Units: ng/uL

ANALYTE	CURVE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,4-Dioxane	Ave	0.4034	0.4205	0.0100	5.21	5.00	4.2	20.0
N-Nitrosodimethylamine	Ave	0.5369	0.5491	0.0100	5.11	5.00	2.3	20.0
Pyridine	Ave	0.9657	0.9797	0.0100	5.07	5.00	1.4	20.0
Methyl methanesulfonate	Ave	0.5820	0.5787	0.0100	4.97	5.00	-0.6	20.0
Benzaldehyde	Qua	0.6652	0.7990	0.0100	4.69	5.00	-6.1	20.0
Aniline	Ave	1.715	1.692	0.0100	4.93	5.00	-1.3	20.0
Phenol	Ave	1.632	1.575	0.8000	4.83	5.00	-3.5	20.0
Bis(2-chloroethyl)ether	Ave	1.106	1.057	0.7000	4.78	5.00	-4.4	20.0
2-Chlorophenol	Ave	1.282	1.261	0.8000	4.92	5.00	-1.6	20.0
1,3-Dichlorobenzene	Ave	1.589	1.567	0.0100	4.93	5.00	-1.4	20.0
1,4-Dichlorobenzene	Ave	1.610	1.569	0.0100	4.87	5.00	-2.6	20.0
Benzyl alcohol	Ave	0.7606	0.6876	0.0100	4.52	5.00	-9.6	20.0
1,2-Dichlorobenzene	Ave	1.532	1.508	0.0100	4.92	5.00	-1.6	20.0
Indene	Ave	2.253	2.183	0.0100	4.84	5.00	-3.1	20.0
2-Methylphenol	Ave	1.132	1.119	0.7000	4.94	5.00	-1.2	20.0
2,2'-oxybis[1-chloropropane]	Ave	1.584	1.549	0.0100	4.89	5.00	-2.2	20.0
N-Nitrosopyrrolidine	Ave	0.5434	0.5051	0.0100	4.65	5.00	-7.0	20.0
Acetophenone	Ave	1.749	1.653	0.0100	4.73	5.00	-5.5	20.0
N-Nitrosodi-n-propylamine	Ave	0.8499	0.8198	0.5000	4.82	5.00	-3.5	20.0
Methylphenol, 3 & 4	Ave	1.223	1.160	0.6000	4.74	5.00	-5.2	20.0
Hexachloroethane	Ave	0.5685	0.5745	0.3000	5.05	5.00	1.1	20.0
Nitrobenzene	Ave	0.3169	0.3349	0.2000	5.28	5.00	5.7	20.0
Isophorone	Ave	0.5497	0.5455	0.4000	4.96	5.00	-0.8	20.0
2-Nitrophenol	Ave	0.1821	0.1902	0.1000	5.22	5.00	4.5	20.0
2,4-Dimethylphenol	Ave	0.3096	0.3114	0.2000	5.03	5.00	0.6	20.0
Benzoic acid	Lin	0.1791	0.1568	0.0100	9.00	10.0	-10.0	20.0
Bis(2-chloroethoxy)methane	Ave	0.3598	0.3503	0.3000	4.87	5.00	-2.6	20.0
2,4-Dichlorophenol	Ave	0.3003	0.3123	0.2000	5.20	5.00	4.0	20.0
1,2,4-Trichlorobenzene	Ave	0.3574	0.3627	0.0100	5.07	5.00	1.5	20.0
Naphthalene	Ave	1.048	1.042	0.7000	4.97	5.00	-0.6	20.0
4-Chloroaniline	Ave	0.4229	0.3959	0.0100	4.68	5.00	-6.4	20.0
2,6-Dichlorophenol	Ave	0.2976	0.3018	0.0100	5.07	5.00	1.4	20.0
Hexachlorobutadiene	Ave	0.2220	0.2343	0.0100	5.28	5.00	5.5	20.0
Caprolactam	Lin	0.0831	0.0773	0.0100	4.33	5.00	-13.3	20.0
4-Chloro-3-methylphenol	Ave	0.2842	0.2695	0.2000	4.74	5.00	-5.2	20.0
2-Methylnaphthalene	Ave	0.7405	0.7054	0.4000	4.76	5.00	-4.7	20.0
1-Methylnaphthalene	Ave	0.6875	0.6614	0.0100	4.81	5.00	-3.8	20.0
Hexachlorocyclopentadiene	Ave	0.3846	0.4078	0.0500	5.30	5.00	6.0	20.0
1,2,4,5-Tetrachlorobenzene	Ave	0.6357	0.7041	0.0100	5.54	5.00	10.7	20.0
2,4,6-Trichlorophenol	Ave	0.3537	0.3971	0.2000	5.61	5.00	12.3	20.0
2,4,5-Trichlorophenol	Ave	0.3801	0.4079	0.2000	5.36	5.00	7.3	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93710/10 Calibration Date: 01/02/2014 12:15

Instrument ID: 732 Calib Start Date: 12/11/2013 05:33

GC Column: Rxi-5SilMS ID: 0.32(mm) Calib End Date: 12/11/2013 08:37

Lab File ID: D01020CC.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,1'-Biphenyl	Ave	1.449	1.478	0.0100	5.10	5.00	2.0	20.0
2-Chloronaphthalene	Ave	1.175	1.281	0.8000	5.45	5.00	9.0	20.0
2-Nitroaniline	Ave	0.2810	0.2963	0.0100	5.27	5.00	5.4	20.0
Dimethyl phthalate	Ave	1.258	1.228	0.0100	4.88	5.00	-2.4	20.0
1,3-Dinitrobenzene	Lin	0.1947	0.1897	0.0100	4.81	5.00	-3.7	20.0
2,6-Dinitrotoluene	Ave	0.2832	0.2783	0.2000	4.91	5.00	-1.7	20.0
Acenaphthylene	Ave	1.760	1.788	0.9000	5.08	5.00	1.6	20.0
3-Nitroaniline	Ave	0.3089	0.2868	0.0100	4.64	5.00	-7.2	20.0
2,4-Dinitrophenol	Lin	0.1825	0.1563	0.0100	9.15	10.0	-8.5	20.0
Acenaphthene	Ave	1.176	1.188	0.9000	5.05	5.00	1.1	20.0
4-Nitrophenol	Ave	0.1479	0.1348	0.0100	9.11	10.0	-8.9	20.0
2,4-Dinitrotoluene	Ave	0.3894	0.3653	0.2000	4.69	5,00	-6.2	20.0
Dibenzofuran	Ave	1.677	1.664	0.8000	4.96	5.00	-0.8	20.0
2,3,5,6-Tetrachlorophenol	Ave	0.3708	0.3442	0.0100	4.64	5.00	-7.2	20.0
2,3,4,6-Tetrachlorophenol	Ave	0.3671	0.3534	0.0100	4.81	5.00	-3.7	20.0
2-Naphthylamine	Ave	0.5439	0.5354	0.0100	4.92	5.00	-1.6	20.0
Diethyl phthalate	Ave	1.209	1.185	0.0100	4.90	5.00	-1.9	20.0
4-Chlorophenyl phenyl ether	Ave	0.7231	0.6981	0.4000	4.83	5.00	-3.5	20.0
4-Nitroaniline	Ave	0.3179	0.2820	0.0100	4.43	5.00	-11.3	20.0
Fluorene	Ave	1.351	1.309	0.9000	4.85	5.00	-3.1	20.0
4,6-Dinitro-2-methylphenol	Lin	0.1380	0.1384	0.0100	10.0	10.0	0.1	20.0
N-Nitrosodiphenylamine	Ave	0.5278	0.5486	0.0100	5.20	5.00	3.9	20.0
1,2-Diphenylhydrazine(as Azobenzene)	Ave	0.6173	0.6988	0.0100	5.66	5.00	13.2	20.0
4-Bromophenyl phenyl ether	Ave	0.2260	0.2393	0.1000	5.29	5.00	5.9	20.0
Hexachlorobenzene	Ave	0.2200	0.2184	0.1000	4.97	5.00	-0.7	20.0
Atrazine	Qua	0.1590	0.1625	0.0100	4.74	5.00	-5.2	20.0
Pentachlorophenol	Lin	0.1688	0.1429	0.0500	9.49	10.0	-5.1	20.0
Phenanthrene	Ave	1.110	1.108	0.7000	4.99	5.00	-0.2	20.0
Anthracene	Ave	1.113	1.116	0.7000	5.01	5.00	0.3	20.0
Carbazole	Ave	1.006	0.9813	0.0100	4.88	5.00	-2.4	20.0
Di-n-butyl phthalate	Ave	1.106	1.097	0.0100	4.96	5.00	-0.8	20.0
Fluoranthene	Ave	1.258	1.202	0.6000	4.77	5.00	-4.5	20.0
Benzidine	Qua	0.2433	0.2024	0.0100	4.72	5.00	-5.6	20.0
Pyrene	Ave	1.081	1.205	0.6000	5.57	5.00	11.4	20.0
Butyl benzyl phthalate	Lin	0.3673	0.4359	0.0100	5.38	5.00	7.5	20.0
3,3'-Dichlorobenzidine	Qua	0.3234	0.3575	0.0100	5.00	5.00	-0.0	20.0
Bis(2-ethylhexyl) phthalate	Lin	0.4963	0.6025	0.0100	5.40	5.00	8.0	20.0
Benzo[a]anthracene	Ave	1.079	1.119	0.8000	5.18	5.00	3.7	20.0
Chrysene	Ave	1.021	1.026	0.7000	5.03	5.00	0.5	20.0
Di-n-octyl phthalate	Qua	0.9594	1.182	0.0100	5.88	5.00	17.5	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93710/10 Calibration Date: 01/02/2014 12:15

Instrument ID: 732 Calib Start Date: 12/11/2013 05:33

Lab File ID: D01020CC.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
7,12-Dimethylbenz(a)anthrace	Ave	0.5682	0.5881	0.0100	5.18	5.00	3.5	20.0
Benzo[b] fluoranthene	Ave	1.232	1.309	0.7000	5.31	5.00	6.3	20.0
Benzo[k] fluoranthene	Ave	1.278	1.237	0.7000	4.84	5.00	-3.2	20.0
Benzo[a]pyrene	Ave	1.095	1.144	0.7000	5.23	5.00	4.5	20.0
<pre>Indeno[1,2,3-cd]pyrene</pre>	Ave	1.160	1.202	0.5000	5.18	5.00	3.6	20.0
Dibenz(a,h)anthracene	Ave	0.9749	1.032	0.4000	5.29	5.00	5.8	20.0
Benzo[g,h,i]perylene	Ave	0.996	1.015	0.5000	5.10	5.00	2.0	20.0
2-Fluorophenol (Surr)	Ave	1.074	1.136		5.29	5.00	5.8	20.0
Phenol-d5 (Surr)	Ave	1.415	1.441		5.09	5.00	1.9	20.0
Nitrobenzene-d5 (Surr)	Ave	0.3040	0.3322	<u>-</u>	5.46	5.00	9.3	20.0
2-Fluorobiphenyl	Ave	1.346	1.385		5.14	5.00	2.9	20.0
2,4,6-Tribromophenol (Surr)	Ave	0.0838	0.0865	0.0100	5.16	5.00	3.2	20.0
Terphenyl-d14 (Surr)	Ave	0.8022	0.8846	<u> </u>	5.51	5.00	10.3	20.0

### FORM V GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

DFTPP Injection Date: 11/07/2013 Lab File ID: S1107DF1.D

Instrument ID: 71 DFTPP Injection Time: 03:02

Analysis Batch No.: 89450

M/E	ION ABUNDANCE CRITERIA	v	ATIVE DANCE
51	30.0 - 60.0 % of mass 198	43.0	
68	Less than 2.0 % of mass 69	0.7	(1.7)1
69	Mass 69 relative abundance	39.9	-
70	Less than 2.0 % of mass 69	0.0	(0.0)1
127	40.0 - 60.0 % of mass 198	46.2	
197	Less than 1.0 % of mass 198	0.0	
198	Base Peak, 100 % relative abundance	100.0	
199	5.0- 9.0 % of mass 198	7.0	
275	10.0 - 30.0 % of mass 198	25.3	
365	Greater than 1.0 % of mass 198	2.7	
441	Present but less than mass 443	10.5	(85.7)3
442	Greater than 40.0 % of mass 198	67.8	
443	17.0 - 23.0 % of mass 442	12.3	(18.1)2

1-Value is % mass 69 2-Value is % mass 442 3-Value is % mass 443

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 180-89450/1	S11070C2.D	11/07/2013	04:37
·	IC 180-89450/2	S11070C3.D	11/07/2013	05:04
	ICIS 180-89450/3	S11070C4.D	11/07/2013	05:30
	IC 180-89450/4	S11070C5.D	11/07/2013	05:57
	IC 180-89450/5	S11070C6.D	11/07/2013	06:23
	IC 180-89450/6	S11070C7.D	11/07/2013	06:50
	IC 180-89450/7	S11070C8.D	11/07/2013	07:16
	IC 180-89450/8	S11070C9.D	11/07/2013	08:11

### GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION FORM VI

Lab Name: TestAmerica Pittsburgh		Job	b No.:	180-28384	384-1				1	Analy B	Batch No	No.: 89	89450	
SDG No.:							!							
Instrument ID: 71	,	25	Column:	Rxi-	5SilMS	ID:	0.32 (mm)	(1)		Heated	Purge:	(X/N)	z	
Calibration Start Date: 11/07/2013	04:37	Cal	libration	End	Date:	11/07/2013	2013	08:11		Calibration	tion ID	): 12367	29	
ANALYTE			RRF			CURVE	COE	COEFFICIENT	#	MIN RRF	%RSD #		R^2 #	
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	TYPE	ш	Ml	M2			%RSD	OR COD	OR COD
2-Methylphenol	1.1974	1.3049	1.3882	1.3912	1.3822	Ave		1.3179		0.7000	5.1	20.0		
Indene	1.3166	2.4825	1.2867	2.4956	2.4829	Ave		2.4250		0.0100	4.0	20.0		!
2,2'-oxybis[1-chloropropane]	2.5713	2.5711	2.6050	2.5117	2.4868	Ave		2.3927		0.0100	9.7	20.0		
N-Nitrosopyrrolidine	0.5623	0.6528	0.6777	0.6944	0.7149	Ave		0.6671		0.0100	7.0	20.0		
Acetophenone	2.0601	2.3114	2.1938	2.1495	2.0878	Ave		2.0559		0.0100	8.1	20.0		
N-Nitrosodi-n-propylamine	1.0566 1.0566	1.1827	1.1820	1.1712	1.1601	Ave		1.0979		0.5000	7.7	20.0		
Methylphenol, 3 & 4	1.3617	1.4204	1.4746	1.4427	1.4330	Ave		1.3809		0.6000	5.1	20.0		
Hexachloroethane	0.5931	0.6546	0.6275	0.6279	0.6326	Ave		0.6238		0.3000	2.9	20.0	:	
Nitrobenzene	0.3318	0.3871	0.3966	0.4025	0.4061	Ave		0.3878	!	0.2000	6.1	20.0		
Isophorone	0.5904	0.6736	0.6608	0.7128	0.7135	Ave		0.6831		0.4000	6.3	20.0		
2-Nitrophenol	+++++	0.1500	0.1610	0.1728	0.1813	Ave		0.1734		0.1000	7.6	20.0		
2,4-Dimethylphenol	0.3082	0.2873	0.3687	0.3655	0.3645	Ave		0.3421		0.2000	8.6	20.0		
Bis(2-chloroethoxy)methane	0.4100	0.4351	0.4380	0.4358	0.4261	Ave		0.4200		0.3000	3.7	20.0		<u></u>
Benzoic acid	+ C	0.0802	0.1076	0.1464	0.1694	Qua	0.4206	5.3915	-0.085	0.0100		20.0	8666.0	0066.0
2,4-Dichlorophenol	0.2176	0.2685	0.2775	0.2831	0.2819	Ave		0.2668		0.2000	7.9	20.0		
1,2,4-Trichlorobenzene	0.3325	0.3344	0.3247	0.3213	0.3090	Ave		0.3144		0.0100	5.1	20.0		
Naphthalene	1.1017	1.0871	1.0923	1.0636	1.0476	Ave		1.0482		0.7000	4.4	20.0		
4-Chloroanlline	0.3941	0.4396	0.4317	0.4452	0.4433	Ave		0.4230		0.0100	4.8	20.0	<u> </u>	
2,6-Dichlorophenol	0.2373	0.2815	0.2778	0.2874	_	Ave	:	0.2695		0.0100	5.9	20.0		1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
Hexachlorobutadiene	0.2074	0.1991	0.1986	0.1957	0.1917	Ave		0.1931		0.0100	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 INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh		Job	b No.:	180-28384	384-1				44	Analy Ba	Batch No.		9450	
SDG No.:														!
Instrument ID: 71		SS	Column:	Rxi	-5SilMS	ID:	0.32 (mm)	( u	14	Heated P	Purge:	(Y/N)	z	
Calibration Start Date: 11/07/2013	04:37	Cal	libration	on End	Date:	11/07/2013	2013	08:11		Calibration	ID	: 1236	57	-
ANALYTE			RRF			CURVE	COE	COEFFICIENT	#	MIN RRF	%RSD #	MAX	R^2 #	MIN R^2
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	TYPE	m m	M1	M2			*KSD	 	OK COD
Caprolactam	++++	0.0862	0.0980	0.1057	0.1050	Ave		0.0958		0.0100	8.4	20.0		
4-Chloro-3-methylphenol	0.0974	0.0917	0.0864	0.3184	0.3169	Ave		0.2970		0.2000	7.3	20.0		
2-Methylnaphthalene	0.3045	0.2995	0.2904	0.7475	0.7239	Ave		0.7143		0.4000	6.1	20.0		
1-Methylnaphthalene	0.6798	0.7148	0.7061	0.6946	0.6732	Ave	_ *	0.6657		0.0100	6.3	20.0		
Hexachlorocyclopentadiene	+ + + + + + + + + + + + + + + + + + + +	0.2778	0.2961	0.3168	0.3245	Ave		0.3232		0.0500	6.8	20.0		
1,2,4,5-Tetrachlorobenzene	0.3416	0.5269	0.3563	0.5325	0.5259	Ave		0.5468		0.0100	4.5	20.0		
2,4,6-Trichlorophenol	0.2345	0.3341	0.3469	0.3458	0.3535	Ave		0.3353		0.2000	12.4	20.0		
2,4,5-Trichlorophenol	0.2532	0.3434	0.3613	0.3742	0.3712	Ave	<del></del>	0.3470		0.2000	11.4	20.0		Ţ——
1,1'-Bipheny1	1.5497	1.5405	1.5237	1.4567	1.4317	Ave		1.4582		0.0100	4.9	20.0		:
2-Chloronaphthalene	1.2396	1.2036	1.1655	1.1564	1.1525	Ave		1.1625		0.8000	3.9	20.0		
2-Nitroaniline	+++++	0.3011	0.3214	0.3572	0.3691	Ave		0.3468	<u> </u>	0.0100	7.3	20.0		
Dimethyl phtholate	1.1508	1.3243	1.3252	1.2755	1.2394	Ave	:	1.2134		0.0100	7.4	20.0		:
1,3-Dinitrobenzene	+++++	0.1409	0.1628	0.1879	0.2008	Ave		0.1853		0.0100	13.1	20.0		
2,6-Dinitrotoluene	+++++	0.2567	0.2799	0.2885	0.2840	Ave		0.2750		0.2000	4.0	20.0	1	V
Acenaphthylene	1.6518	1.8563	1.8804	1.8570	1.8109	Ave	L	1.7813		0.9000	4.7	20.0		- <del></del>
3-Nitroaniline	+++++	0.2849	0.2853	0.3119	0.3136	Ave		0.2949		0.0100	4.3	20.0		
2,4-Dinitrophenol	+++++	+++++	0.0450	0.0736	0.0997	Ona	0.8543	8.1984 -(	-0.434	0.0100		20.0	9666.0	0.9900
Acenaphthene	1.2253	1.2277	1.2051	1.1886	1.1367	Ave		1,1451		0.9000	6.7	20.0		<u> </u>
4-Nitrophenol	+++++	0.1145	0.1335	0.1584	0.1655	Ave		0.1503		0.0100	12.6	20.0		
Z,4-Dinitrotoluene	0.1455	0.3201	0.3356	0.3601	0.3575	Qua	0.000.0	2.7907 0.	0.1181	0.2000		20.0	7666.0	0.9900

Note: The ml coefficient is the same as Ave RRF for an Ave curve type.

### FORM VI GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Pittsburgh		Ϋ́	Job No.:	180-28384	384-1			!		Analy Ba	Batch No.:		89450	
SDG No.:								!						i
Instrument ID: 71	-	gc	Column:	Rxi-	5SilMS	ID: 0	0.32 (mm)			Heated E	Purge:	(Y/N)	Z	
Calibration Start Date: 11/07/2013	04:37	3   	Calibration	on End	Date:	11/07/2013		08:11		Calibration	ion ID	: 12367	29	
ANALYTE			RRF			CURVE	COEF	COEFFICIENT	**=	MIN RRF	%RSD #	MAX	R^2	# MIN R^2
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	TYPE	В	M1	M2			% X U	OR COD	00 00 00 00 00 00 00 00 00 00 00 00 00
Dibenzofuran	1.7824	1.7454		1.6811	1.6111	Ave		1.6361		0.008.0	7.3	20.0		
2,3,5,6-Tetrachlorophenol	1.5459	0.2785	1.4897	0.3167	0.3193	Ave	0	0.3065		0.0100	4.5	20.0		
2,3,4,6-Tetrachlorophenol	++++	0.3226	0.3198	0.3394	0.3319	Ave	0	0.3145	:	0.0100	0.9	20.0		
2-Naphthylamine	0.5152	1.0989	0.6180	0.5868	0.5069	Qua –(	-0.147	1.7694 0	0.3728	0.0100	i	20.0	0.9984	0066.0
Diethyl phthalate	1.1998	1.3684	1.3578	1.3112	1.2525	Ave		1,2187		0.0100	10.2	20.0	<u>:</u>	-
Hexadecane	0.5593	0.6380	0.6666	0.6560	0.6441	Ave	0	0.6083			8.2			
4-Chlorophenyl phenyl ether	0.7012	0.5878	0.7039	0.6863	0.6486	Ave	0	0.6554		0.4000	8.6	20.0		
4-Nitroaniline	+ + 0	0.2669	0.2819	0.3008	0.2838	Ave	0	0.2710		0.0100	7.0	20.0		ļ
Fluorene	1.3985	1.4293	1.4128	1.3547	1.2864	Ave	1	1.2925		0.9000	9.6	20.0		
4,6-Dinitro-2-methylphenol	+ + + + + + + + + + + + + + + + + + + +	0.0487	0.0612	0.0879	0.1048	O eno	0.4473 8	8.7273	-0.324	0.0100		20.0	0.9998	0.9900
N-Nitrosodiphenylamine	0.5043	0.5398	0.5549	0.5593	0.5509	Ave	0	0.5458		0.0100	e	20.0		
1,2-Diphenylhydrazine (as Azobenzene)	0.7220	0.8387	0.8689	0.8809	0.8920	Ave		0.8706		0.0100	7.7	20.0		
4-Bromophenyl phenyl ether	0.2141	0.2279	0.2323	0.2308	0.2287	Ave		0.2246		0.1000	3.0	20.0		
Hexachlorobenzene	0.2524	0.2427	0.2415	0.2360	0.2302	Ave	-	0.2285		0.1000	7.9	20.0	-	
Atrazine	+++++	0.1986	0.1571	0.1597	0.1576	Ave	0	0.1595		0.0100	11.4	20.0		
Pentachlorophenol	0.0338	0.1047	0.1026	0.1236	0.1302	Qua	0.0740 7	7.8434 0	0.1390	0.0500		20.0	9666.0	0066.0
n-Octadecane	2.0219	2.8412	2.9579	2.9477	2.9585	Ave	2	2.6291			13.7			
Phenanthrene	1.1943	1.1342	1.1324	1.1086	1.0810	Ave		1.1061		0.7000	4.2	20.0		
Anthracene	0.9944	137	1.1393	1.1265	1.0898	Ave		1.0667		0.7000	0.9	20.0		-
Carbazole	0.8481	0.9952	0.9982	0.9862	0.9514 Ave	Ave	0	0.9248		0.0100	7.0	20.0		

Note: The ml coefficient is the same as Ave RRF for an Ave curve type.

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## Page 1115 of 1943

## GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION FORM VI

		dot	on q	180-28384-1	384-1					Analy Batch No.:	tch No.		89450	
SDG No.: Instrument ID: 71		25	Column:		Rxi-5SilMS	ID: 0	.32 (mm)			Heated P	Purge: (	(Y/N)	z	
Calibration Start Date: 11/07/2013	04:37	Cal	libration	End	Date:	11/07/201	<u></u>	08:11		Calibration	ion ID:	12367	57	
ANALYTE			RRF			CURVE	COEF	COEFFICIENT	! ##	MIN RRE	%RSD #		R^2 #	MIN
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	TYPE		M1	M2			*RSD	OR COD	OR COD
Di-n-butyl phthalate	0.7721	1.1150	1.1259	1.1953	1.1984	Ave		6260.	-	0.0100	12.4	20.0		
Fluoranthene	0.9881	1.1456	1.1230	1.1487	1.1035	Ave		1.0652		0.6000	6.9	20.0		·
Benzidine	++++	++++	0.1233	0.1303	0.1330	Qua 0.	0.3231 5	5.9257	-0.641	0.0100		20.02	0.9978	0.9900
Pyrene	1.2648	1.2909	1.3521	1.3583	1.3615	Ave		1.3136		0.009.0	3.0	20.0	:	
Butyl benzyl phthalate	+ CT + C	0.4092	0.4325	0.4879	0.5236	Ave	0	0.4955		0.0100	11.1	20.0	-	
3,3'-Dichlorobenzidine	+++++	0.2816	0.2893	0.3209	0.3458	Ave	.0	0.3426		0.0100	13.6	20.0		
Bis(2-ethylhexyl) phthalate	) + c	0.4879	0.5454	0.6174	0.6865	Ave	0	0.6393		0.0100	14.5	20.0		· - <u> </u>
Benzo[a]anthracene	1.0032	1.1014	1.1257	1,1188	1.1262	Ave		1.0901		0.8000	3.7	20.0		
Chrysene	0.9924	1.0222	1.0424	1.0320	1.0134	Ave		1.0040		0.7000	2.7	20.0		
Di-n-octyl phthalate	+++++	1.1643	1.0349	1.1745	1.3341	Ave	1	1.2164		0.0100	8.4	20.0	:	
7,12-Dimethylbenz(a)anthracene	0.4466	0.5043	0.6109	0.6184	0.6315	Ave	0	0.5520		0.0100	12.4	20.0		
Benzo[b]fluoranthene	1.1139	1.2848	1.3735	1.3372	1.3541	Ave		1.2510		0.7000	8.5	20.0		
Benzo[k]fluoranthene	1.0597	1.2775	1.2805	1.2933	1.3153	Ave	-	1.2145	i	0.7000	8.1	20.0		·
Benzo[a]pyrene	0.8256	1.0204	1.0733	1.1246	1.1763	Ave	1	1.0726		0.7000	10.4	20.0		:
Indeno[1,2,3-cd]pyrene	1.0437	1.1012	1.1804	1.2227	1.3149	Ave		1.2385		0.5000	6.6	20.0		
Dibenz(a,h)anthracene	0.8549	0.9410	1.0234	1.0463	1.1300	Ave		1.0466		0.4000	10.0	20.0		
Benzo[g,h,1]perylene	0.8929	0.9688	1.0014	1.0264	1.1138	Ave	-	1.0607		0.5000	8.6	20.0	<u> </u>	
2-Fluorophenol (Surr)	1.2628	1.1927	1.2462	1.2480	1.2659	Ave		1.2281			4.5	20.0		
Phenol-d5 (Surr)	1.5249	1.6624	1.7385	1.7357	1.7469	Ave		1.6796			4.4	20.0		
Nitrobenzene-d5 (Surr)	0.3281	0.3891	0.3835	0.3864	0.3926	Ave	0	0.3765			5.6	20.0		
							-			i				

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

## GC/MS SEMI VOA INITIAL CALIBRATION DATA INTERNAL STANDARD CURVE EVALUATION FORM VI

# MIN R^2

MAX R^2

MIN RRF %RSD #

====

COEFFICIENT

11/07/2013 08:11

ID: 0.32 (mm)

GC Column: Rxi-5SilMS

Job No.: 180-28384-1

TestAmerica Pittsburgh

Lab Name: SDG No.: Calibration End Date:

04:37

11/07/2013

Calibration Start Date:

Instrument ID:

ANALYTE

 $M_2$ 

Ξ

щ

LVL 5

LVL 4

LVL 3

RRF

CURVE

Calibration ID: 12367

Heated Purge: (Y/N)

89450

Analy Batch No.:

0.9900

9866.0

20.0 20.02

0.0100

10.178 1.2422

0.0421

Qua

1.2991 Ave 0.0952

1.3299

1,3635 .2850

1.4066 LVL 2 LVL 7

1.3692 LVL 1 LVL 6

1.2548

0.9428

0.9848 Ave

0.9848 0.0922

0.0859 0.0886 0.9909 0.8950

0.0891 0.9492 0.9145

1.2698 0.0346 0.0933 0.8864 0.9364

2,4,6-Tribromophenol (Surr)

2-Fluorobiphenyl

Terphenyl-d14 (Surr)

4.1

20.0

4.4

### FORM V GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

DFTPP Injection Date: 01/01/2014 Lab File ID: S0101DF1.D

Instrument ID: 71 DFTPP Injection Time: 04:01

Analysis Batch No.: 93752

M/E	ION ABUNDANCE CRITERIA	,	JATIVE DANCE
51	30.0 - 60.0 % of mass 198	33.5	
68	Less than 2.0 % of mass 69	0.0	(0.0)1
69	Mass 69 relative abundance	36.1	
70	Less than 2.0 % of mass 69	0.3	(0.7)1
127	40.0 - 60.0 % of mass 198	44.4	
197	Less than 1.0 % of mass 198	0.0	
198	Base Peak, 100 % relative abundance	100.0	
199	5.0- 9.0 % of mass 198	6.8	
275	10.0 - 30.0 % of mass 198	25.9	
365	Greater than 1.0 % of mass 198	3.0	
441	Present but less than mass 443	12.3	(92.7)3
442	Greater than 40.0 % of mass 198	68.4	
443	17.0 - 23.0 % of mass 442	13.3	(19.5)2

1-Value is % mass 69 2-Value is % mass 442 3-Value is % mass 443

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 180-93752/22	S01010C2.D	01/01/2014	04:44
	MB 180-93531/1-A	S0101001.D	01/01/2014	05:41
	LCS 180-93531/2-A	S0101002.D	01/01/2014	06:35
SD-185-0-1 MS	180-28384-23 MS	S0101003.D	01/01/2014	08:22
SD-185-0-1 MSD	180-28384-23 MSD	S0101004.D	01/01/2014	08:48
SD-183-0-1	180-28384-21	S0101005.D	01/01/2014	09:15
SD-184-0-2	180-28384-22	S0101006.D	01/01/2014	09:42
SD-185-0-1	180-28384-23	S0101007.D	01/01/2014	10:08
SD-186-0-1	180-28384-24	S0101008.D	01/01/2014	10:35
SD-187-0-1	180-28384-25	S0101009.D	01/01/2014	11:01
SD-188-0-1	180-28384-26	S0101010.D	01/01/2014	11:28
SD-189-0-1	180-28384-27	S0101011.D	01/01/2014	11:55
SD-190-0-1	180-28384-28	S0101012.D	01/01/2014	12:22
SD-191-0-1	180-28384-29	S0101013.D	01/01/2014	12:48
SD-192-0-1	180-28384-30	S0101014.D	01/01/2014	13:15
SD-193-0-1	180-28384-31	S0101015.D	01/01/2014	13:42
SD-194-0-1	180-28384-32	S0101016.D	01/01/2014	14:09

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93752/22 Calibration Date: 01/01/2014 04:44

Instrument ID: 71 Calib Start Date: 11/07/2013 04:37

GC Column: Rxi-5SilMS ID: 0.32(mm) Calib End Date: 11/07/2013 08:11

Lab File ID: S01010C2.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	용D	MAX %D
1,4-Dioxane	Ave	0.4681	0.3553	0.0100	3.80	5.00	-24.1*	20.0
N-Nitrosodimethylamine	Ave	0.6583	0.4986	0.0100	3.79	5.00	-24.3*	20.0
Pyridine	Ave	1.147	0.8659	0.0100	3.77	5.00	-24.5*	20.0
Methyl methanesulfonate	Ave	0.7540	0.5888	0.0100	3.90	5.00	-21.9*	20.0
Benzaldehyde	Qua	0.8334	0.8710	0.0100	4.58	5.00	-8.3	20.0
Aniline	Ave	2.130	1.812	0.0100	4.25	5.00	-14.9	20.0
Phenol	Ave	1.912	1.723	0.8000	4.51	5.00	-9.9	20.0
Bis(2-chloroethyl)ether	Ave	1.431	1.221	0.7000	4.27	5.00	-14.7	20.0
2-Chlorophenol	Ave	1.409	1.363	0.8000	4.84	5.00	-3.2	20.0
1,3-Dichlorobenzene	Ave	1.528	1.566	0.0100	5.12	5.00	2.4	20.0
1,4-Dichlorobenzene	Ave	1.535	1.601	0.0100	5.21	5.00	4.3	20.0
Benzyl alcohol	Ave	0.9074	0.9032	0.0100	4.98	5.00	-0.5	20.0
1,2-Dichlorobenzene	Ave	1.486	1.564	0.0100	5.26	5.00	5.2	20.0
Indene	Ave	2.425	. 2.405	0.0100	4.96	5.00	-0.8	20.0
2-Methylphenol	Ave	1.318	1.307	0.7000	4.96	5.00	-0.8	20.0
2,2'-oxybis[1-chloropropane]	Ave	2.393	1.574	0.0100	3.29	5.00	-34.2*	20.0
N-Nitrosopyrrolidine	Ave	0.6671	0.6391	0.0100	4.79	5.00	-4.2	20.0
Acetophenone	Ave	2.056	1.991	0.0100	4.84	5.00	-3.1	20.0
N-Nitrosodi-n-propylamine	Ave	1.098	1.028	0.5000	4.68	5.00	-6.4	20.0
Methylphenol, 3 & 4	Ave	1.381	1,433	0.6000	5.19	5.00	3.8	20.0
Hexachloroethane	Ave	0.6238	0.6066	0.3000	4.86	5.00	-2.8	20.0
Nitrobenzene	Ave	0.3878	0.3313	0.2000	4.27	5,00	-14.6	20.0
Isophorone	Ave	0.6831	0.6072	0.4000	4.44	5.00	-11.1	20.0
2-Nitrophenol	Ave	0.1734	0.1745	0.1000	5.03	5.00	0.6	20.0
2,4-Dimethylphenol	Ave	0.3421	0.3320	0.2000	4.85	5.00	-3.0	20.0
Benzoic acid	Qua	0.1528	0.2050	0.0100	12.6	10.0	26.5*	20.0
Bis(2-chloroethoxy)methane	Ave	0.4200	0.3720	0.3000	4.43	5.00	-11.4	20.0
2,4-Dichlorophenol	Ave	0.2668	0.2918	0.2000	5.47	5.00	9.4	20.0
1,2,4-Trichlorobenzene	Ave	0.3144	0.3283	0.0100	5.22	5.00	4.4	20.0
Naphthalene	Ave	1.048	1.049	0.7000	5.00	5.00	0.0	20.0
4-Chloroaniline	Ave	0.4230	0.4053	0.0100	4.79	5.00	-4.2	20.0
2,6-Dichlorophenol	Ave	0.2695	0.2984	0.0100	5.54	5.00	10.7	20.0
Hexachlorobutadiene	Ave	0,1931	0.2068	0.0100	5.35	5.00	7.1	20.0
Caprolactam	Ave	0.0958	0.1200	0.0100	6.26	5,00	25.3*	20.0
4-Chloro-3-methylphenol	Ave	0,2970	0,3265	0.2000	5.50	5.00	9.9	20.0
2-Methylnaphthalene	Ave	0.7143	0.7817	0.4000	5.47	5.00	9.4	20.0
1-Methylnaphthalene	Ave	0.6657	0.7284	0.0100	5.47	5.00	9.4	20.0
Hexachlorocyclopentadiene	Ave	0.3232	0.3277	0.0500	5.07	5.00	1.4	20.0
1,2,4,5-Tetrachlorobenzene	Ave	0.5468	0.5397	0,0100	4.93	5.00	-1.3	20.0
2,4,6-Trichlorophenol	Ave	0.3353	0.3490	0.2000	5.21	5.00	4.1	20.0
2,4,5-Trichlorophenol	Ave	0.3333	0.3752	0.2000	5.41	5.00	8.1	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93752/22 Calibration Date: 01/01/2014 04:44

Instrument ID: 71 Calib Start Date: 11/07/2013 04:37

GC Column: Rxi-5SilMS ID: 0.32(mm) Calib End Date: 11/07/2013 08:11

Lab File ID: S01010C2.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	왕D	MAX %D
1,1'-Biphenyl	Ave	1.458	1.413	0.0100	4.84	5.00	-3.1	20.0
2-Chloronaphthalene	Ave	1.162	1.161	0.8000	4.99	5.00	-0.1	20.0
2-Nitroaniline	Ave	0.3468	0.3144	0.0100	4.53	5.00	-9.4	20.0
Dimethyl phthalate	Ave	1.213	1.286	0.0100	5.30	5.00	5.9	20.0
1,3-Dinitrobenzene	Ave	0.1853	0.2104	0.0100	5.68	5.00	13.6	20.0
2,6-Dinitrotoluene	Ave	0.2750	0.3018	0.2000	5.49	5.00	9.7	20.0
Acenaphthylene	Ave	1.781	1.823	0.9000	5.12	5.00	2.3	20.0
3-Nitroaniline	Ave	0.2949	0.3280	0.0100	5.56	5.00	11.2	20.0
Acenaphthene	Ave	1.145	1.170	0.9000	5.11	5.00	2.2	20.0
2,4-Dinitrophenol	Qua	0.1005	0.1145	0.0100	12.7	10.0	26.6*	20.0
4-Nitrophenol	Ave	0.1503	0.1758	0.0100	11.7	10.0	17.0	20.0
2,4-Dinitrotoluene	Qua	0.3096	0.4179	0.2000	5.99	5.00	19.8	20.0
Dibenzofuran	Ave	1.636	1.708	0.8000	5.22	5.00	4.4	20.0
2,3,5,6-Tetrachlorophenol	. Ave	0.3065	0.3441	0.0100	5.61	5.00	12.3	20.0
2,3,4,6-Tetrachlorophenol	Ave	0.3145	0.3725	0.0100	5.92	5.00	18.4	20.0
2-Naphthylamine	Qua	0.5527	0.4670	0.0100	4.05	5.00	-18.9	20.0
Diethyl phthalate	Ave	1.219	1.325	0.0100	5.44	5.00	8.7	20.0
4-Chlorophenyl phenyl ether	Ave	0.6554	0.7116	0.4000	5.43	5.00	8.6	20.0
Fluorene	Ave	1.292	1.438	0.9000	5.56	5.00	11.2	20.0
4-Nitroaniline	Ave	0.2710	0.3636	0.0100	6.71	5.00	34.2*	20.0
4,6-Dinitro-2-methylphenol	Qua	0.0944	0.1208	0.0100	12.2	10.0	22.1*	20.0
N-Nitrosodiphenylamine	Ave	0.5458	0.5299	0.0100	4.85	5.00	-2.9	20.0
1,2-Diphenylhydrazine(as Azobenzene)	Ave	0.8706	0.7144	0.0100	4.10	5.00	-17.9	20.0
4-Bromophenyl phenyl ether	Ave	0.2246	0.2197	0.1000	4.89	5.00	-2.2	20.0
Hexachlorobenzene	Ave	0.2285	0.2262	0.1000	4.95	5.00	-1.0	20.0
Atrazine	Ave	0.1595	0.1714	0.0100	5.37	5.00	7.4	20.0
Pentachlorophenol	Qua	0.1078	0.1286	0.0500	10.4	10.0	4.4	20.0
Phenanthrene	Ave	1.106	1.132	0.7000	5.12	5.00	2.3	20.0
Anthracene	Ave	1.067	1.175	0.7000	5.51	5.00	10.1	20.0
Carbazole	Ave	0.9248	1.101	0.0100	5.95	5.00	19.0	20.0
Di-n-butyl phthalate	Ave	1.098	1.250	0.0100	5.69	5.00	13.8	20.0
Fluoranthene	Ave	1.065	1.296	0.6000	6.08	5.00	21.6*	20.0
Benzidine	Qua	0.1610	0.1009	0.0100	4.24	5.00	-15.2	20.0
Pyrene	Ave	1.314	1.086	0.6000	4.13	5.00	-17.4	20.0
Butyl benzyl phthalate	Ave	0.4955	0.4693	0.0100	4.74	5.00	-5.3	20.0
3,3'-Dichlorobenzidine	Ave	0.3426	0.4034	0.0100	5.89	5.00	17.7	20.0
Bis(2-ethylhexyl) phthalate	Ave	0.6393	0.6723	0.0100	5.26	5.00	5.2	20.0
Benzo[a]anthracene	Ave	1.090	1.117	0.8000	5.12	5.00	2.5	20.0
Chrysene	Ave	1.004	1.027	0.7000	5.12	5.00	2.3	20.0
Di-n-octyl phthalate	Ave	1.216	1.158	0.0100	4.76	5.00	-4.8	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

Lab Sample ID: CCVIS 180-93752/22 Calibration Date: 01/01/2014 04:44

Calib Start Date: 11/07/2013 04:37

Instrument ID: 71

GC Column: Rxi-5SilMS ID: 0.32(mm) Calib End Date: 11/07/2013 08:11

Lab File ID: S01010C2.D

Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
7,12-Dimethylbenz(a)anthrace	Ave	0.5520	0.5418	0.0100	4.91	5.00	-1.8	20.0
Benzo[b]fluoranthene	Ave	1.251	1.207	0.7000	4.83	5.00	-3.5	20.0
Benzo[k]fluoranthene	Ave	1.214	1.171	0.7000	4.82	5.00	-3.6	20.0
Benzo[a]pyrene	Ave	1.073	1.111	0.7000	5.18	5.00	3.6	20.0
Indeno[1,2,3-cd]pyrene	Ave	1.238	1.332	0.5000	5.38	5.00	7.6	20.0
Dibenz(a,h)anthracene	Ave	1.047	1.136	0.4000	5.43	5.00	8.5	20.0
Benzo[g,h,i]perylene	Ave	1.061	1.118	0.5000	5.27	5.00	5.4	20.0
2-Fluorophenol (Surr)	Ave	1.228	1.054		4.29	5.00	-14.1	20.0
Phenol-d5 (Surr)	Ave	1.680	1.527		4.55	5.00	-9.1	20.0
Nitrobenzene-d5 (Surr)	Ave	0.3765	0.3293		4.37	5.00	-12.5	20.0
2-Fluorobipheny1	Ave	1.322	1.275		4.82	5.00	-3.5	20.0
2,4,6~Tribromophenol (Surr)	Qua	0.0823	0.0945	0.0100	5.05	5.00	0.9	20.0
Terphenyl-d14 (Surr)	Ave	0.9428	0.8074		4.28	5.00	-14.4	20.0

#### FORM V

### GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab File ID: S0104DF1.D DFTPP Injection Date: 01/04/2014

Instrument ID: 71 DFTPP Injection Time: 07:39

Analysis Batch No.: 93785

M/E	ION ABUNDANCE CRITERIA		ATIVE DANCE
51	30.0 - 60.0 % of mass 198	34.3	
68	Less than 2.0 % of mass 69	0.0	(0.0)1
69	Mass 69 relative abundance	36.3	
70	Less than 2.0 % of mass 69	0.4	(1.2)1
127	40.0 - 60.0 % of mass 198	44.8	
197	Less than 1.0 % of mass 198	0.0	
198	Base Peak, 100 % relative abundance	100.0	
199	5.0- 9.0 % of mass 198	6.8	
275	10.0 - 30.0 % of mass 198	26.1	
365	Greater than 1.0 % of mass 198	3.1	
441	Present but less than mass 443	12.8	(97.5)3
442	Greater than 40.0 % of mass 198	66.5	
443	17.0 - 23.0 % of mass 442	13.1	(19.7)2

1-Value is % mass 69 2-Value is % mass 442 3-Value is % mass 443

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 180-93785/3	S01040C1.D	01/04/2014	07:53
SD-190-0-1 DL	180-28384-28 DL	S0104012.D	01/04/2014	13:38

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93785/3 Calibration Date: 01/04/2014 07:53

Instrument ID: 71 Calib Start Date: 11/07/2013 04:37

GC Column: Rxi-5SilMS ID: 0.32(mm) Calib End Date: 11/07/2013 08:11

Lab File ID: S01040C1.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,4-Dioxane	Ave	0.4681	0.3889	0.0100	4.15	5.00	-16.9	20.0
N-Nitrosodimethylamine	Ave	0.6583	0.5153	0.0100	3.91	5.00	-21.7*	20.0
Pyridine	Ave	1.147	0.8860	0.0100	3.86	5.00	-22.8*	20.0
Methyl methanesulfonate	Ave	0.7540	0.5431	0.0100	3.60	5.00	-28.0*	20.0
Benzaldehyde	Qua	0.8334	0.6088	0.0100	3.36	5.00	-32.9*	20.0
Aniline	Ave	2.130	1.568	0.0100	3.68	5.00	-26.4*	20.0
Phenol	Ave	1.912	1.538	0.8000	4.02	5.00	-19.5	20.0
Bis(2-chloroethyl)ether	Ave	1.431	1.112	0.7000	3.88	5.00	-22.3*	20.0
2-Chlorophenol	Ave	1.409	1.307	0.8000	4.64	5.00	-7.2	20.0
1,3-Dichlorobenzene	Ave	1.528	1.544	0.0100	5.05	5.00	1.0	20.0
1,4-Dichlorobenzene	Ave	1.535	1.581	0.0100	5.15	5.00	3.0	20.0
Benzyl alcohol	Ave	0.9074	0.7853	0.0100	4.33	5.00	-13.4	20.0
1,2-Dichlorobenzene	Ave	1.486	1.532	0.0100	5.15	5.00	3.1	20.0
Indene	Ave	2.425	2.250	0.0100	4.64	5.00	-7.2	20.0
2-Methylphenol	Ave	1.318	1.163	0.7000	4.41	5.00	-11.7	20.0
2,2'-oxybis[1-chloropropane]	Ave	2.393	1.396	0.0100	2.92	5.00	-41.7*	20.0
N-Nitrosopyrrolidine	Ave	0.6671	0.5096	0.0100	3.82	5.00	-23.6*	20.0
Acetophenone	Ave	2.056	1.729	0.0100	4.20	5.00	-15.9	20.0
N-Nitrosodi-n-propylamine	Ave	1.098	0.8540	0.5000	3.89	5.00	-22.2*	20.0
Methylphenol, 3 & 4	Ave	1.381	1.230	0.6000	4.45	5.00	-10.9	20.0
Hexachloroethane	Ave	0.6238	0.5841	0.3000	4.68	5.00	-6.4	20.0
Nitrobenzene	Ave	0.3878	0.3361	0.2000	4.33	5.00	-13.3	20.0
Isophorone	Ave	0.6831	0.5716	0.4000	4.18	5.00	-16.3	20.0
2-Nitrophenol	Ave	0.1734	0.1745	0.1000	5.03	5.00	0.6	20.0
2,4-Dimethylphenol	Ave	0,3421	0.3208	0.2000	4.69	5.00	-6.2	20.0
Benzoic acid	Qua	0.1528	0.1457	0.0100	9.49	10.0	-5.1	20.0
Bis(2-chloroethoxy)methane	Ave	0.4200	0.3533	0.3000	4.21	5.00	-15.9	20.0
2,4-Dichlorophenol	Ave	0.2668	0.2975	0.2000	5.57	5.00	11.5	20.0
1,2,4-Trichlorobenzene	Ave	0.3144	0.3445	0.0100	5.48	5.00	9.6	20.0
Naphthalene	Ave	1.048	1.055	0.7000	5.03	5.00	0.6	20.0
4-Chloroaniline	Ave	0.4230	0.3871	0.0100	4.58	5.00	-8.5	20.0
2,6-Dichlorophenol	Ave	0.2695	0.2972	0.0100	5.51	5.00	10.3	20.0
Hexachlorobutadiene	Ave	0.1931	0.2282	0.0100	5.91	5.00	18.2	20.0
Caprolactam	Ave	0.0958	0.0901	0.0100	4.70	5.00	-5.9	20.0
4-Chloro-3-methylphenol	Ave	0.2970	0.2847	0.2000	4.79	5.00	-4.2	20.0
2-Methylnaphthalene	Ave	0.7143	0.7009	0.4000	4.91	5.00	-1.9	20.0
1-Methylnaphthalene	Ave	0.6657	0.6592	0.0100	4.95	5.00	-1.0	20.0
Hexachlorocyclopentadiene	Ave	0.3232	0.3333	0.0500	5.16	5.00	3.1	20.0
1,2,4,5-Tetrachlorobenzene	Ave	0.5468	0.5831	0.0100	5.33	5.00	6.6	20.0
2,4,6-Trichlorophenol	Ave	0.3353	0.3565	0.2000	5.32	5.00	6.3	20.0
2,4,5-Trichlorophenol	Ave	0.3470	0.3721	0.2000	5.36	5.00	7.2	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93785/3 Calibration Date: 01/04/2014 07:53

Instrument ID: 71 Calib Start Date: 11/07/2013 04:37

GC Column: Rxi-5SilMS ID: 0.32(mm) Calib End Date: 11/07/2013 08:11

Lab File ID: S01040C1.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,1'-Biphenyl	Ave	1.458	1.400	0.0100	4.80	5.00	-4.0	20.0
2-Chloronaphthalene	Ave	1.162	1.140	0.8000	4.90	5.00	-2.0	20.0
2-Nitroaniline	Ave	0.3468	0.2889	0.0100	4.16	5.00	-16.7	20.0
Dimethyl phthalate	Ave	1.213	1.210	0.0100	4.99	5.00	-0.3	20.0
1,3-Dinitrobenzene	Ave	0.1853	0.1922	0.0100	5.19	5.00	3.7	20.0
2,6-Dinitrotoluene	Ave	0.2750	0.2818	0.2000	5.12	5.00	2.5	20.0
Acenaphthylene	Ave	1.781	1.856	0.9000	5.21	5.00	4.2	20.0
3-Nitroaniline	Ave	0.2949	0.3160	0.0100	5.36	5.00	7.2	20.0
Acenaphthene	Ave	1.145	1.177	0.9000	5.14	5.00	2.8	20.0
2,4-Dinitrophenol	Qua	0.1005	0.0763	0.0100	9.61	10.0	-3.9	20.0
4-Nitrophenol	Ave	0.1503	0.1635	0.0100	10.9	10.0	8.8	20.0
2,4-Dinitrotoluene	Qua	0.3096	0.4044	0.2000	5.79	5.00	15.8	20.0
Dibenzofuran	Ave	1.636	1.792	0.8000	5.48	5.00	9.5	20.0
2,3,5,6-Tetrachlorophenol	Ave	0.3065	0.3219	0.0100	5.25	5.00	5.0	20.0
2,3,4,6-Tetrachlorophenol	Ave	0.3145	0.3499	0.0100	5.56	5.00	11.3	20.0
2-Naphthylamine	Qua	0.5527	0.5186	0.0100	4.63	5.00	-7.5	20.0
Diethyl phthalate	Ave	1.219	1.352	0.0100	5.55	5.00	10.9	20.0
4-Chlorophenyl phenyl ether	Ave	0.6554	0.7342	0.4000	5.60	5.00	12.0	20.0
4-Nitroaniline	Ave	0.2710	0.3412	0.0100	6.29	5.00	25.9*	20.0
Fluorene	Ave	1.292	1.487	0.9000	5.75	5.00	15.0	20.0
4,6-Dinitro-2-methylphenol	Qua	0.0944	0.1001	0.0100	10.4	10.0	4.4	20.0
N-Nitrosodiphenylamine	Ave	0.5458	0.5423	0.0100	4.97	5.00	-0.6	20.0
1,2-Diphenylhydrazine(as Azobenzene)	Ave	0.8706	0.7409	0.0100	4.26	5.00	-14.9	20.0
4-Bromophenyl phenyl ether	Ave	0.2246	0.2286	0.1000	5.09	5.00	1.8	20.0
Hexachlorobenzene	Ave	0.2285	0.2295	0.1000	5.02	5.00	0.4	20.0
Atrazine	Ave	0.1595	0.1661	0.0100	5.21	5.00	4.1	20.0
Pentachlorophenol	Qua	0.1078	0.1082	0.0500	8.82	10.0	-11.8	20.0
Phenanthrene	Ave	1.106	1.144	0.7000	5.17	5.00	3.4	20.0
Anthracene	Ave	1.067	1.167	0.7000	5.47	5.00	9.4	20.0
Carbazole	Ave	0.9248	1.057	0.0100	5.72	5.00	14.3	20.0
Di-n-butyl phthalate	Ave	1.098	1.251	0.0100	5.70	5.00	13.9	20.0
Fluoranthene	Ave	1.065	1.257	0.6000	5.90	5.00	18.0	20.0
Benzidine	Qua	0.1610	0.1872	0.0100	6.70	5.00	34.0*	20.0
Pyrene	Ave	1.314	1.145	0.6000	4.36	5.00	-12.8	20.0
Butyl benzyl phthalate	Ave	0.4955	0.4916	0.0100	4.96	5.00	-0.8	20.0
3,3'-Dichlorobenzidine	Ave	0.3426	0.3905	0.0100	5.70	5.00	14.0	20.0
Bis(2-ethylhexyl) phthalate	Ave	0.6393	0.6956	0.0100	5.44	5.00	8.8	20.0
Benzo[a]anthracene	Ave	1.090	1.133	0.8000	5.20	5.00	3.9	20.0
Chrysene	Ave	1.004	1.062	0.7000	5.29	5.00	5.8	20.0
Di-n-octyl phthalate	Ave	1.216	1.313	0.0100	5.40	5.00	7.9	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCVIS 180-93785/3 Calibration Date: 01/04/2014 07:53

Instrument ID: 71 Calib Start Date: 11/07/2013 04:37

Lab File ID: S01040C1.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
7,12-Dimethylbenz(a)anthrace	Ave	0.5520	0.5641	0.0100	5.11	5.00	2.2	20.0
Benzo[b] fluoranthene	Ave	1.251	1.270	0.7000	5.08	5.00	1.6	20.0
Benzo[k]fluoranthene	Ave	1.214	1.228	0.7000	5.06	5.00	1.1	20.0
Benzo[a]pyrene	Ave	1.073	1.127	0.7000	5.25	5.00	5.1	20.0
Indeno[1,2,3-cd]pyrene	Ave	1.238	1.387	0.5000	5.60	5.00	12.0	20.0
Dibenz(a,h)anthracene	Ave	1.047	1.165	0.4000	5.57	5.00	11.3	20.0
Benzo[g,h,i]perylene	Ave	1.061	1.196	0.5000	5.64	5.00	12.8	20.0
2-Fluorophenol (Surr)	Ave	1.228	1.085		4.42	5.00	-11.7	20.0
Phenol-d5 (Surr)	Ave	1.680	1.382		4.11	5.00	-17.7	20.0
Nitrobenzene-d5 (Surr)	Ave	0.3765	0.3292		4.37	5.00	-12.6	20.0
2-Fluorobiphenyl	Ave	1.322	1.305		4.94	5.00	-1.3	20.0
2,4,6-Tribromophenol (Surr)	Qua	0.0823	0.0940	0.0100	5.02	5.00	0.4	20.0
Terphenyl-d14 (Surr)	Ave	0.9428	0.8380		4.44	5.00	-11.1	20.0

### FORM II GC SEMI VOA SURROGATE RECOVERY

Lab	Name:	TestAmerica	Pittsburgh	Job No.:	180-28384-1

SDG No.:

Matrix: Sediment Level: Low

GC Column (2): RTX-1701 ID: 0.53 (mm)

Client Sample ID	Lab Sample ID	TCX2 #	DCB2 #
SD-163-0-1	180-28384-1	102	103
SD-164-0-1	180-28384-2	104	180 2
SD-165-0-1	180-28384-3	99	105
SD-166-0-1	180-28384-4	88	73
SD-167-0-1	180-28384-5	69	69
SD-168-0-1	180-28384-6	102	131
SD-169-0-1	180-28384-7	76	74
SD-170-0-1	180-28384-8	71	64
SD-171-0-1	180-28384-9	0 X I	D 0 X I
SD-172-0-3	180~28384-10	0 X I	0 X I
SD-173-0-3	180-28384-11	0 X I	D 0 X I
SD-174-0-1	180-28384-12	96	131
SD-175-0-1	180-28384-13	120	D 180 D 2
SD-176-0-3	180-28384-14	89	88
SD-177-0-2	180-28384-15	93	84
SD-178-0-1	180-28384-16	75	62
SD-179-0-1	180-28384-17	94	81
SD-180-0-1	180-28384-18	115	61
SD-181-0-1	180-28384-19	0 X I	D 0 X I
SD-182-0-3	180-28384-20	0 X I	D 0 X I
SD-183-0-1	180-28384-21	0 X I	D 0 X I
SD-184-0-2	180-28384-22	89	105
SD-185-0-1	180-28384-23	123	103
SD-186-0-1	180-28384-24	87	64
SD-187-0-1	180-28384-25	86	65
SD-188-0-1	180-28384-26	77	64
SD-189-0-1	180-28384-27	82	70
SD-190-0-1	180-28384-28	81	53
SD-191-0-1	180-28384-29	82	89
SD-192-0-1	180-28384-30	65	54
SD-193-0-1	180-28384-31	78	68
SD-194-0-1	180-28384-32	91	72
	MB 180-93257/1-C	92	82
	MB 180-93333/1-C	76	70
	<u> </u>	<u> </u>	

QC LIMITS

30-150

20-150

TCX = Tetrachloro-m-xylene
DCB = DCB Decachlorobiphenyl (Surr)

# Column to be used to flag recovery values

### FORM II GC SEMI VOA SURROGATE RECOVERY

Lab Name	: TestAmerica	Pittsburgh	Job No.	: 180-28384-1	
SDG No.:					
Matrix:	Sediment		Level:	Low	

GC Column (2): RTX-1701 ID: 0.53 (mm)

Client Sample ID	Lab Sample ID	TCX2 #	DCB2 #
-	LCS	91	81
	180-93257/2-C		
	LCS	84	74
	180-93333/2-C		
SD-165-0-1 MS	180-28384-3 MS	92	95
SD-185-0-1 MS	180-28384-23 MS	103	47
SD-165-0-1 MSD	180-28384-3 MSD	88	94
SD-185-0-1 MSD	180-28384-23 MSD	98	65

QC LIMITS

TCX = Tetrachloro-m-xylene
DCB = DCB Decachlorobiphenyl (Surr)

30-150 20-150

# Column to be used to flag recovery values

FORM II 8082A

### FORM IV GC SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Pitts	burgh	Job No.: 180-28384-1
SDG No.:		
Lab Sample ID: MB 180-9325	7/1-C	
Matrix: Sediment		Date Extracted: 12/26/2013 03:40
Lab File ID: (1)		Lab File ID:(2) T1230582.D
Date Analyzed:(1)	·	Date Analyzed:(2) 12/27/2013 21:15
<pre>Instrument ID:(1)</pre>		Instrument ID:(2) GC10
GC Column:(1)	ID:	GC Column:(2) RTX-1701 ID: 0.53(mm)

#### THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

GLIEUE GIMBLE TO		DATE	DATE
CLIENT SAMPLE ID	LAB SAMPLE ID	ANALYZED 1	ANALYZED 2
SD-163-0-1	180-28384-1		12/27/2013 04:27
SD-164-0-1	180-28384-2		12/27/2013 04:58
SD-165-0-1	180-28384-3		12/27/2013 05:30
SD-165-0-1 MS	180-28384-3 MS		12/27/2013 06:01
SD-165-0-1 MSD	180-28384-3 MSD		12/27/2013 06:33
SD-166-0-1	180-28384-4		12/27/2013 07:04
SD-167-0-1	180-28384-5		12/27/2013 07:36
SD-168-0-1	180-28384-6		12/27/2013 08:07
SD-169-0-1	180-28384-7		12/27/2013 08:39
SD-170-0-1	180-28384-8		12/27/2013 09:10
SD-174-0-1	180-28384-12		12/27/2013 14:26
SD-176-0-3	180-28384-14		12/27/2013 15:28
SD-177-0-2	180-28384-15		12/27/2013 16:00
SD-178-0-1	180-28384-16		12/27/2013 16:31
SD-179-0-1	180-28384-17		12/27/2013 17:03
SD-171-0-1	180-28384-9		12/27/2013 19:09
SD-172-0-3	180-28384-10		12/27/2013 19:40
SD-173-0-3	180-28384-11		12/27/2013 20:12
	LCS 180-93257/2-C		12/27/2013 21:46
SD-175-0-1	180-28384-13		12/28/2013 08:16
SD-180-0-1	180-28384-18		12/28/2013 08:48
SD-181-0-1	180-28384-19		12/28/2013 11:58
SD-182-0-3	180-28384-20		12/28/2013 12:29

Page 1 of 2

Lab Name: TestAmerica Pittsburgh

Job No.: 180-28384-1

SDG No.:

93257 Batch Number:

Batch Start Date: 12/26/13 03:20

Batch Analyst: Geehring, Kevin

Batch End Date: 12/26/13 09:47 Batch Method: 3541

op-p/pcb sur 00010	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL	25 uL
GCMATRIXWORKS 00009		25 uL	25 uL	25 uL	:														-		
InitialAmount	30.0 g	30.0 g	30.1 g	30.1 g	30.1 g	30.0 g	30.1 g	30.0 g	30.0 g	30.0 g	30.1 g	30.1 g	30.0 g	30.0 g	30.0 g	30.1 g	30.0 g	30.0 g	30.2 g	30.1 g	30.0 g
FinalAmount	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL	1.0 mL
Basis			 [ <del>-</del> ]	£	T	F	E	<u> </u>	E	Ŀ	EH	E	E-i	E	F	E	E	-	E	E+	E
Method Chain	3541, 3665A, 3660B, 8082A	3541, 3665A, 3660B, 8082A	$\sim$	3541, 3665A,	3541, 3665A,	3541, 3665A,	3541, 3665A, 3660B, 8082A	3541, 3665A, 3660B, 8082A	3541, 3665A, 3660B, 8082A		3541, 3665A,		3541, 3665A,				3541, 3665A,	3541, 3665A,	3541, 3665A,	m	m
Client Sample ID	-		SD-165-0-1	SD-165-0-1	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-167-0-1	SD-168-0-1	SD-169-0-1	SD-170-0-1	SD-171-0-1	SD-172-0-3	SD-173-0-3	SD-174-0-1	SD-175-0-1	SD-176-0-3	SD-177-0-2	SD-178-0-1	SD-179-0-1
Lab Sample ID	MB 180-93257/1	LCS 180-93257/2	180-28384-A-3	180-28384-A-3 MSD	180-28384-A-1	180-28384-A-2	180-28384-A-3	180-28384-A-4	180-28384-A-5	180-28384-A-6	180-28384-A-7	180-28384-A-8	180-28384-A-9	180-28384-A-10	180-28384-A-11	180-28384-A-12	180-28384-A-13	180-28384-A-14	180-28384-A-15	180-28384-A-16	180-28384-A-17

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.

8082A

Page 2 of 2

# GC SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Pittsburgh

Job No.: 180-28384-1

SDG No.:

Batch Number: 93257

Batch Start Date: 12/26/13 03:20

Batch Analyst: Geehring, Kevin

Batch Method: 3541

Batch End Date: 12/26/13 09:47

ıı			
op-p/pcb sur 00010	25 uL	25 ul	25 uL
GCMATRIXWORKS 00009			
InitialAmount	30.0 g	.30.1 g	30.0 g
FinalAmount	1.0 mL	1.0 mL	1.0 mL
Basis	E	E+	E
Method Chain	3541, 3665A, 3660B, 8082A	3541, 3665A, 3660B, 8082A	3541, 3665A,
Lab Sample ID Client Sample ID Method Chain Basis		SD-181-0-1	SD-182-0-3
Lab Sample ID	180-28384-A-18   SD-180-0-1	180-28384-A-19   SD-181-0-1	180-28384-A-20 SD-182-0-3

	Batch Notes
Balance ID	1120122641
Batch Comment	sox # 1 - 2 - 3 - 4
Person's name who did the concentration	kg
Exchange Solvent Lot #	1017776
Exchange Solvent Name	Hexane
Magnesium Sulfate Lot #	1055012
N-evap #	
Na2SO4 Lot Number	1047333
Person's name who did the prep	KG KG
Solvent	Hexane/acetone
Solvent Lot #	1055182
Uncorrected N-evap Temperature	32 Degrees C

Description		
Basis		Total/NA
asis	-	E

### FORM I GC SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1 SDG No.: Client Sample ID: Lab Sample ID: MB 180-93257/1-C Lab File ID: T1230582.D Matrix: Sediment Analysis Method: 8082A Date Collected: Extraction Method: 3541 Date Extracted: 12/26/2013 03:40 Sample wt/vol: 30.0(g) Date Analyzed: 12/27/2013 21:15 Con. Extract Vol.: 1.0(mL) Dilution Factor: 1 GC Column: RTX-1701 ID: 0.53(mm) Injection Volume: 1(uL) GPC Cleanup:(Y/N) N % Moisture: Analysis Batch No.: 93552 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT Q	RL	MDL
12674-11-2	PCB-1016	ND	0.42	0.062
11104-28-2	PCB-1221	ND	0.42	0.080
11141-16-5	PCB-1232	ND	0.42	0.071
53469-21-9	PCB-1242	ND	0.42	0.068
12672-29-6	PCB-1248	ND	0.42	0.039
11097-69-1	PCB-1254	ND	0.42	0.059
11096-82-5	PCB-1260	ND	0.42	0.059

CAS NO.	SURROGATE	%REC	Q	LIMITS
2051-24-3	DCB Decachlorobiphenyl (Surr)	82		20-150
877-09-8	Tetrachloro-m-xylene	92		30-150

### FORM III GC SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name	: TestAmerica Pitts	sburgh	Job No.: 180-28384-1				
SDG No.:							
Matrix:	Sediment	Level: Low	Lab File ID: 1	[1230583.D			
Lab ID:	LCS 180-93257/2-C	·	Client ID:				

	SPIKE ADDED	LCS CONCENTRATION	LCS %	QC LIMITS	#
COMPOUND	(ug/Kg)	(ug/Kg)	REC	REC	
PCB-1016	33.3	29.5	88	50-120	
PCB-1260	33.3	28.6	86	50-120	

 $<sup>\</sup>mbox{\#}$  Column to be used to flag recovery and RPD values FORM III  $8082\mbox{\ensuremath{\text{A}}}$ 

### FORM III GC SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Pitt	sburgh	Job No.: 1	80-28384-1			
SDG No.:				-		
Matrix: Sediment	Level: Low	Lab File II	T1230553.D			
Lab ID: 180-28384-3 MS		Client ID:	SD-165-0-1 MS			
	SPIKE	SAMPLE	MS	MS	oc	
	ADDED	CONCENTRATION	CONCENTRATION	ક	LIMITS	#
COMPOUND	(ug/Kg)	(ug/Kg)	(ug/Kg)	ŖEC	REC	
PCB-1016	43.6	ND	47.4	109	50-120	
PCB-1260	43.6	99	117	40)	50-120	F

 $<sup>\</sup>mbox{\#}$  Column to be used to flag recovery and RPD values FORM III  $8082\mbox{\ensuremath{\text{R}}}$ 

### FORM III GC SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Pitt	sburgh	Job No.: 1	80-283	884-1			
SDG No.:							
Matrix: Sediment	Level: Low	Lab File II	): <u>T1</u>	230554	.D		
Lab ID: 180-28384-3 MSD		Client ID:	SD-1	55-0 <b>-</b> 1	MSD		
	SPIKE	MSD	MSD		QC LI	MITS	
	ADDED	CONCENTRATION	olo Olo	용			#
COMPOUND	(ug/Kg)	(ug/Kg)	REC	RPD	RPD	REC	
PCB-1016	43.6	46.3	<u>10</u> 6	2	30	50-120	
PCB-1260	43.6	113	(3)3	3	30	50-120	F

 $<sup>\</sup>mbox{\#}$  Column to be used to flag recovery and RPD values FORM III 8082A

### FORM IV GC SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Pi	ttsburgh	Job No.: 180-28384-1
SDG No.:		
Lab Sample ID: MB 180-93	3333/1-C	
Matrix: Sediment		Date Extracted: 12/27/2013 02:40
Lab File ID:(1)		Lab File ID:(2) T1230600.D
Date Analyzed: (1)		Date Analyzed:(2) 12/28/2013 06:42
Instrument ID:(1)		Instrument ID:(2) GC10
GC Column:(1)	ID:	GC Column:(2) RTX-1701 ID: 0.53(mm)

### THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

		DATE	DATE
CLIENT SAMPLE ID	LAB SAMPLE ID	ANALYZED 1	ANALYZED 2
SD-184-0-2	180-28384-22		12/27/2013 23:21
SD-185-0-1	180-28384-23		12/27/2013 23:52
SD-185-0-1 MS	180-28384-23 MS		12/28/2013 00:24
SD-185-0-1 MSD	180-28384-23 MSD		12/28/2013 00:55
SD-186-0-1	180-28384-24		12/28/2013 01:27
SD-187-0-1	180-28384-25		12/28/2013 01:58
SD-188-0-1	180-28384-26		12/28/2013 02:30
SD-189-0-1	180-28384-27		12/28/2013 03:01
SD-190-0-1	180-28384-28		12/28/2013 03:33
SD-191-0-1	180-28384-29		12/28/2013 04:04
SD-192-0-1	180-28384-30		12/28/2013 04:36
SD-193-0-1	180-28384-31		12/28/2013 05:07
SD-194-0-1	180-28384-32		12/28/2013 05:39
	LCS 180-93333/2-C		12/28/2013 07:13
SD-183-0-1	180-28384-21		12/28/2013 10:23

2

Page 1 of

## GC SEMI VOA BATCH WORKSHEET

02:40 12/27/13 Job No.: 180-28384-1 Batch Start Date: TestAmerica Pittsburgh 93333 Batch Number: Lab Name: SDG No.:

Batch End Date: 12/27/13 08:37

3541

Batch Method:

Geehring, Kevin

Batch Analyst:

sur op-p/pcb : 25 uL 25 uL 25 uL 25 uL 25 uL 25 uL αŢ 25 uL 25 uL 25 uL 25 uL μŢ μĽ 25 uL  $_{\rm uL}$ ηŢ 25 25 25 25 25 GCMATRIXWORKS 00009 25 uL 25 uL ηŢ 25 InitialAmount 30.1 g 30.09 30.0g 30.2 g 30.09 30.0g 30.0g 30.0g 30.1 g 30.1 g 30.0g 30.0g 30.1 g 30.0 g 30.1 g FinalAmount 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL 1.0 mL Basis Method Chain 3541, 3665A, 3660B, 8082A 3541, 3665A, 3660B, 8082A 3541, 36654, 3560B, 8082A 3541, 36634, 3560B, 8082A 3660B, 8082A 3541, 3665A, 3560B, 8082A 3541, 3665A, 3560B, 8082A 3561, 3665A, 3541, 3665A, 3660B, 8082A, 366 3541, 3665A, 3660B, 8082A, 3541, 3665A, 3660B, 8082A 3660B, 8082A 3541, 3665A, 3660B, 8082A ΩĪ Client Sample SD-185-0-1 SD-185-0-1 SD-183-0-1 SD-184-0-2 SD-185-0-1 SD-186-0-1 SD-187-0-1 SD-188-0-1 SD-189-0-1 SD-190-0-1 SD-191-0-1 SD-192-0-1 SD-193-0-1 SD-194-0-1 Lab Sample ID LCS 180-93333/2 180-28384-A-23 180-28384-A-23 180-28384-A-23 180-28384-A-24 180-28384-A-26 180-28384-A-29 180-28384-A-30 180-28384-A-32 180-28384-A-21 180-28384-A-22 180-28384-A-25 180-28384-A-27 180-28384-A-28 180-28384-A-31 MB 180-93333/1

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.

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Page 2 of 2

### GC SEMI VOA BATCH WORKSHEET

Job No.: 180-28384-1 Lab Name: TestAmerica Pittsburgh

SDG No.:

Batch Start Date: 93333 Batch Number:

3541 Batch Method:

02:40 12/27/13

Geehring, Kevin

Batch Analyst:

Batch End Date: 12/27/13 08:37

Bato	Batch Notes
Balance ID	1120122641
Batch Comment	sox # 2 - 3 - 4
Person's name who did the concentration	kg
Exchange Solvent Lot #	1017776
Exchange Solvent Name	Hexane
Magnesium Sulfate Lot #	1055012
N-evap #	. 2
Na2SO4 Lot Number	1047333
Person's name who did the prep	kg kg
Solvent	Hexane/acetone
Solvent Lot #	1055182
Uncorrected N-evap Temperature	32 Degrees C

Basis Description	Total/NA	
Basis	T To	

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 I GC SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1 SDG No.: Lab Sample ID: MB 180-93333/1-C Client Sample ID: Matrix: Sediment Lab File ID: T1230600.D Analysis Method: 8082A Date Collected: Extraction Method: 3541 Date Extracted: 12/27/2013 02:40 Sample wt/vol: 30.0(g) Date Analyzed: 12/28/2013 06:42 Con. Extract Vol.: 1.0(mL) Dilution Factor: 1 GC Column: RTX-1701 ID: 0.53 (mm) Injection Volume: 1(uL) GPC Cleanup: (Y/N) N % Moisture: Analysis Batch No.: 93552 Units: ug/Kg,

CAS NO.	COMPOUND NAME	RESULT	Q RL	MDL
12674-11-2	PCB-1016	ND ·	0.42	0.062
11104-28-2	PCB-1221 .	ND	0.42	0.080
11141-16-5	PCB-1232	ND	0.42	0.071
53469-21-9	PCB-1242	ND	0.42	0.068
12672-29-6	PCB-1248	ND	0.42	0.039
11097-69-1	PCB-1254	ND	0.42	0.059
11096-82-5	PCB-1260	ND	0.42	0.059

CAS NO.	SURROGATE	%REC	Q	:	LIMITS
2051-24-3	DCB Decachlorobiphenyl (Surr)	70			20-150
877-09 <b>-</b> 8	Tetrachloro-m-xylene	76			30-150

### FORM III GC SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name	: TestAmerica Pitt	sburgh	Job N	o.: <u>180-28384-1</u>	<u> </u>		
SDG No.:							
Matrix:	Sediment	Level: Low	Lab F	ile ID: T12306	01.D		
Lab ID:	LCS 180-93333/2-C		Clien	t ID:			
		SP	IKE	LCS	LCS	QC	
		AD	DED	CONCENTRATION	8	LIMITS	#
	COMPOUND	(ug	/Kg)	(ug/Kg)	REC	REC	
PCB-101	16		33.3	26	5.3 79	50-120	
PCB-126	50	<del></del>	33.3	25	5.5 77	50-120	

 $<sup>\</sup>mbox{\#}$  Column to be used to flag recovery and RPD values FORM III  $\mbox{8082A}$ 

### FORM III GC SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Matrix: Sediment Level: Low Lab File ID: <u>T1230588.D</u>

Lab ID: 180-28384-23 MS Client ID: SD-185-0-1 MS

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC	QC LIMITS REC	#
PCB-1016	44.7	ND	67.0	(150	50-120	F
PCB-1260	44.7	550	579	66	50-120	4

 $<sup>\</sup>mbox{\#}$  Column to be used to flag recovery and RPD values FORM III  $8082\mbox{\ensuremath{A}}$ 

### FORM III GC SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Pitt	sburgh	Job No.: 180	0-28384	4-1		
SDG No.:						
Matrix: Sediment	Level: Low	Lab File ID:	T123	0589.D		
Lab ID: 180-28384-23 MSD	· .	Client ID: S	SD-185	-0-1 MSD		
	SPIKE	MSD	MSD	OC L	IMITS	
	ADDED	CONCENTRATION	ક	% · ·-		# .
COMPOUND	(ug/Kg)	(ug/Kg) F	REC F	RPD RPD	REC	:
PCB-1016	44.5	. 58.8	(132)	13 30	50-120	
PCB-1260	44.5	553	(8)	5 30	50-120	(4)

 $<sup>\</sup>mbox{\#}$  Column to be used to flag recovery and RPD values FORM III  $\mbox{8082A}$ 

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## GC SEMI VOA BATCH WORKSHEET

Lab Name: Tes	Lab Name: TestAmerica Pittsburgh	Job No.: 180-28384-1	1-1		ı			
SDG No.:								
Batch Number:	93283	Batch Start Date: 12/26/13 10:24	12/26/13	13 10:24	Batch Analyst: Gupta, Ashok	Gupta,	Ashok	
Batch Method:	3665A	Batch End Date:						

ount																					
itialAmount FinalAmount	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	2 mL 2 mL	
Method Chain Basis Ini	3665A, 3660B, 8082A			3660B, 8082A T T 3665A, 3660B			İ	3660B, 8082A	3665A, T		3660B, 8082A i 3665A, T	İ	-		3665A, TT		3665A, 9082A T				3660B 8082A
Client Sample ID M	36	36	SD-165-0-1 36	36 SD-165-0-1 36	SD-163-0-1	SD-164-0-1 36	SD-165-0-1	SD-166-0-1	SD-167-0-1 36	SD-168-0-1 36	36 SD-169-0-1 36	SD-170-0-1 36	SD-171-0-1 36	SD-172-0-3	SD-173-0-3	SD-174-0-1	SD-175-0-1 36	SD-176-0-3 36	SD-177-0-2	SD-178-0-1	200
Lab Sample ID	MB 180-93257/1-A	LCS 180-93257/2-A	0-28384-A-3-B	MS 180-28384-A-3-C MSD	-28384-A-1-D	180-28384-A-2-B	180-28384-A-3-D	180-28384-A-4-B	180-28384-A-5-B	180-28384-A-6-B	180-28384-A-7-B	180-28384-A-8-B	180-28384-A-9-B	180-28384-A-10-	80-28384-A-11-	80-28384-A-12-	80-28384-A-13-	80-28384-A-14-	80-28384-A-15-	80-28384-A-16-	

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.

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### GC SEMI VOA BATCH WORKSHEET

Job No.: 180-28384-1

Lab Name: TestAmerica Pittsburgh

SDG No.:							.		
Batch Number:	93283		Ba	Batch Start Date:	12/26/13 10:24	0:24	Batch Analyst: Gupta, Ashok	Gupta, Ashok	!
Batch Method: 3665A	3665A		Ва	Batch End Date:					
Lab Sample ID	Lab Sample ID Client Sample ID Method Chain Basis	Method Chain	Basis	InitialAmount	FinalAmount				
180-28384-A-18-  SD-180-0-1	SD-180-0-1	3665A,		2 шТ	2 mL				
180-28384-A-19-   SD-181-0-1	SD-181-0-1	3665A, 8082A	₽	2 m.L	2 mL				
180-28384-A-20- B	SD-182-0-3	3665A, 3660B, 8082A	· : : : :	2 mL	2 mL				

Batch Notes

Basis Description	Total/NA
Basis	T. To

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.

8082A

Page 1 of 2

Lab Name: TestAmerica Pittsburgh

Job No.: 180-28384-1

SDG No.:

93284 Batch Number:

12/26/13 Batch Start Date:

10:25

Batch Analyst: Gupta, Ashok

Batch End Date: Batch Method: 3660B

				77																				
GCTBASOLUTION 00015	2 m.L	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL		2 mL										
FinalAmount	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 mL	2 m.L	2 mL						
InitialAmount	2 mL	2 mL	2 mL	2 mL	2 mL	2 m.L	2 mL																	
Basis			L	E	H	L	T	T	F	E+	H	T	H	H	E	H	H	E	E→	E	F	H	l H	E
Method Chain	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A	3660B, 8082A
Client Sample ID			SD-165-0-1	SD-165-0-1	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-167-0-1	SD-168-0-1	SD-169-0-1	SD-170-0-1	SD-171-0-1	SD-172-0-3	SD-173-0-3	SD-174-0-1	SD-175-0-1	SD-176-0-3	SD-177-0-2	SD-178-0-1	SD-179-0-1	SD-180-0-1	SD-181-0-1	SD-182-0-3
Lab Sample ID	MB 180-93257/1-B	LCS 180-93257/2-B	180-28384-A-3-E	180-28384-A-3-F MSD	180-28384-A-1-E	180-28384-A-2-C	180-28384-A-3-G	180-28384-A-4-C	180-28384-A-5-C	180-28384-A-6-C	180-28384-A-7-C	180-28384-A-8-C	180-28384-A-9-C	180-28384-A-10-	180-28384-A-11-	180-28384-A-12-	180-28384-A-13-	180-28384-A-14-	180-28384-A-15-	180-28384-A-16-	180-28384-A-17-	180-28384-A-18-	180-28384-A-19-	180-28384-A-20- C

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.

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## GC SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Pittsburgh	Job No.: 180-28384-1		
SDG No.:			
Satch Number: 93284	Batch Start Date: 12/26/13 10:25	Batch Analyst: Gupta, Ashok	Gupta, Ashok
Satch Method: 3660B	Batch End Date:	-	
Ba	Batch Notes		
Batch Comment	H2S04 lot-52267 is used for cleanup ref-3665A	r	
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.

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Page 1 of 1

Lab Name: Tes	TestAmerica Pittsburgh	nrgh	dot	b No.: 180-28384-1	84-1			
SDG No.:								
Batch Number:	93363		Ba	Batch Start Date:	: 12/27/13 10:34	Batch Analyst:	Gupta, Ashok	
Batch Method:	3665A		Ва	Batch End Date:				
Lab Sample ID	Client Sample ID	Method Chain	Basis	InitialAmount	FinalAmount			
MB 180-93333/1-A		3665A, 3660B, 8082A		2 mL	2 mL			
LCS 180-93333/2-A				2 m.L	2 ш.Г			
180-28384-A-23- A MS	SD-185-0-1		E-I	2 mL	2 mL			
180-28384-A-23-	SD-185-0-1		L	2 mL	2 mL			
B MSD 180-28384-A-21-	SD-183-0-1		L	2 mL	2 mL			
A 180-28384-A-22-	SD-184-0-2		E	2 mL	2 mL			
A 180-28384-A-23-	SD-185-0-1		L	2 mL	2 mL			
C 180-28384-A-24-	SD-186-0-1	3660B, 8082A 3665A,	: 1	2 mL	2 mL			
A 180-28384-A-25-	SD-187-0-1	3660B, 8082A 3665A,	E	2 m.L	2 mL			
A   180-28384-A-26-	SD-188-0-1	3660B, 8082A 3665A,	E	2 m.L	2 mL			
A 180-28384-A-27-	SD-189-0-1	3660B, 8082A						:
A 180-28384-A-28-	SD-190-0-1		E					
180-28384-A-29-	SD-191-0-1	3665A, 8082A	E	2 mL	2 mL			
180-28384-A-30-	SD-192-0-1	1	H	2 mL	2 mL			
180-28384-A-31-	SD-193-0-1	1	E	2 mL	2 mL			
180-28384-A-32- A	SD-194-0-1		E+	2 mL	2 mL			
		Bat	Batch Notes	Ø				
Basis E	Basis Description							

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.

Page 1 of 1

## GC SEMI VOA BATCH WORKSHEET

Lab Name: Tes	TestAmerica Pittsburgh	urgh	Job	ob No.: 180-28384-1	34-1			
SDG No.:	:							
Batch Number:	93364		Ba†	Batch Start Date:	12/27/13	10:35	Batch Analyst: Gupta, Ashok	
Batch Method:	3660B		Ba†	atch End Date:				
Lab Sample ID	Client Sample ID	Method Chain	Basis	InitialAmount	FinalAmount	GCTBASOLUTION 00015		
MB 180-93333/1-B		3660B, 8082A		2 mL	2 mL	2 mL		
LCS 180-93333/2-B		3660B, 8082A		2 mL	2 mL	2 mL		
180-28384-A-23-	SD-185-0-1	3660B, 8082A	E	2 mL	2 mL	2 mL		
1180-28384-A-23-	SD-185-0-1	3660B, 8082A	₽	2 mL	2 mL			
180-28384-A-21-	SD-183-0-1	3660B, 8082A		2 mL	2 mL	2 mL		
180-28384-A-22-	SD-184-0-2	3660B, 8082A	E	2 mL	2 mL	2 mL		
180-28384-A-23-	SD-185-0-1	3660B, 8082A	E⊣	2 mL	2 mL	2 mL		
180-28384-A-24-	SD-186-0-1	3660B, 8082A	T	2 mL	2 mL	2 mL		
B 180-28384-A-25-	SD-187-0-1	3660B, 8082A	T	2 mL	2 mL	2 mL		
180-28384-A-26-	SD-188-0-1	3660B, 8082A	E	2 mL	2 mL	2 mL		
180-28384-A-27-	SD-189-0-1	3660B, 8082A	. 🗗	2 mL ·	2 mL	2 mL		
180-28384-A-28-	SD-190-0-1	3660B, 8082A	 E-	2 mL	2 mL	2 mL		
180-28384-A-29-	SD-191-0-1	3660B, 8082A		2 mL	2 mL	2 mL		
180-28384-A-30-	SD-192-0-1	3660B, 8082A	F	2 mL	2 mL	2 mL		
180-28384-A-31-	SD-193-0-1	3660B, 8082A	L	2 mL	2 mL	2 mL		
B 180-28384-A-32- B	SD-194-0-1	3660B, 8082A	·	2 mL	2 mL	2 mL		
		Bat	Batch Notes					
Batch Comment			H2S04	1 lot-53267 is used	for cleanup	ref-3665A		
	Rasis 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 SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Instrument ID: GC10 Start Date: 12/26/2013 17:57

Analysis Batch Number: 93552 End Date: 12/28/2013 15:07

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
IC 180-93552/1		12/26/2013 17:57	1	T1230530.D	RTX-1701 0.53(mm)
IC 180-93552/2		12/26/2013 18:28	1	T1230531.D	RTX-1701 0.53(mm)
IC 180-93552/3		12/26/2013 19:00	1	T1230532.D	RTX-1701 0.53 (mm)
IC 180-93552/4		12/26/2013 19:31	1	T1230533.D	RTX-1701 0.53(mm)
IC 180-93552/5		12/26/2013 20:03	1		RTX-1701 0.53 (mm)
IC 180-93552/6		12/26/2013 20:34	1		RTX-1701 0.53 (mm)
IC 180-93552/7		12/26/2013 21:06	1	T1230536.D	RTX-1701 0.53 (mm)
IC 180-93552/8	:	12/26/2013 21:37		T1230537.D	RTX-1701 0.53 (mm)
IC 180-93552/9		12/26/2013 22:09	1	T1230538.D	RTX-1701 0.53(mm)
ICRT 180-93552/10		12/26/2013 22:40	_ 1	T1230539.D	RTX-1701 0.53 (mm)
IC 180-93552/11		12/26/2013 23:12	1	T1230540.D	RTX-1701 0.53 (mm)
IC 180-93552/12		12/26/2013 23:43	1	T1230541.D	RTX-1701 0.53(mm)
IC 180-93552/13		12/27/2013 00:15	1	T1230542.D	RTX-1701 0.53(mm)
ICV 180-93552/14		12/27/2013 00:46	1	T1230543.D	RTX-1701 0.53(mm)
ICV 180-93552/15		12/27/2013 01:18	1	T1230544.D	RTX-1701 0.53(mm)
ICV 180-93552/16		12/27/2013 01:49	1	T1230545.D	RTX-1701 0.53 (mm)
ICV 180-93552/17		12/27/2013 02:21	1	T1230546.D	RTX-1701 0.53(mm)
ICV 180-93552/18		12/27/2013 02:52	1		RTX-1701 0.53 (mm)
ICV 180-93552/19		12/27/2013 03:24	1		RTX-1701 0.53(mm)
ICV 180-93552/20		12/27/2013 03:55	1	T1230549.D	RTX-1701 0.53(mm)
180-28384-1	SD-163-0-1	12/27/2013 04:27	5	T1230550.D	RTX-1701 0.53 (mm)
180-28384-2	SD-164-0-1	12/27/2013 04:58	5	T1230551.D	RTX-1701 0.53(mm)
180-28384-3	SD-165-0-1	12/27/2013 05:30	5	T1230552.D	RTX-1701 0.53(mm)
180-28384-3 MS	SD-165-0-1 MS	12/27/2013 06:01	5	T1230553.D	RTX-1701 0.53(mm)
180-28384-3 MSD	SD-165-0-1 MSD	12/27/2013 06:33	5	T1230554.D	RTX-1701 0.53(mm)
180-28384-4	SD-166-0-1	12/27/2013 07:04	5	T1230555.D	RTX-1701 0.53(mm)
180-28384-5	SD-167-0-1	12/27/2013 07:36	5	T1230556.D	RTX-1701 0.53(mm)
180-28384-6	SD-168-0-1	12/27/2013 08:07	5	T1230557.D	RTX-1701 0.53(mm)
180-28384-7	SD-169-0-1	12/27/2013 08:39	5	T1230558.D	RTX-1701 0.53(mm)
180-28384-8	SD-170-0-1	12/27/2013 09:10	5	T1230559.D	RTX-1701 0.53(mm)
CCV 180-93552/31		12/27/2013 13:22	1	T1230567.D	RTX-1701 0.53(mm)
180-28384-12	SD-174-0-1	12/27/2013 14:26	5	T1230569.D	RTX-1701 0.53(mm)
180-28384-14	SD-176-0-3	12/27/2013 15:28	5	T1230571.D	RTX-1701 0.53(mm)
180-28384-15	SD-177-0-2	12/27/2013 16:00	5	T1230572.D	RTX-1701 0.53(mm)
180-28384-16	SD-178-0-1	12/27/2013 16:31	5	T1230573.D	RTX-1701 0.53(mm)
180-28384-17	SD-179-0-1	12/27/2013 17:03	5	T1230574.D	RTX-1701 0.53(mm)
180-28384-9	SD-171-0-1	12/27/2013 19:09	1000	T1230578.D	RTX-1701 0.53(mm)
180-28384-10	SD-172-0-3	12/27/2013 19:40	1000	T1230579.D	RTX-1701 0.53 (mm)
180-28384-11	SD-173-0-3	12/27/2013 20:12	5000	T1230580.D	RTX-1701 0.53(mm)
MB 180-93257/1-C		12/27/2013 21:15	1	T1230582.D	RTX-1701 0.53(mm)
LCS 180-93257/2-C		12/27/2013 21:46	1	T1230583.D	RTX-1701 0.53(mm)
CCV 180-93552/42		12/27/2013 22:18		T1230584.D	RTX-1701 0.53(mm)
180-28384-22	SD-184-0-2	12/27/2013 23:21	10	T1230586.D	RTX-1701 0.53(mm)
180-28384-23	SD-185-0-1	12/27/2013 23:52	10	T1230587.D	RTX-1701 0.53(mm)
180-28384-23 MS	SD-185-0-1 MS	12/28/2013 00:24	10	T1230588.D	RTX-1701 0.53 (mm)

### GC SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Instrument ID: GC10 Start Date: 12/26/2013 17:57

Analysis Batch Number: 93552 End Date: 12/28/2013 15:07

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION	LAB FILE ID	COLUMN ID
180-28384-23 MSD	SD-185-0-1 MSD	12/28/2013 00:55	10	T1230589.D	RTX-1701 0.53(mm)
180-28384-24	SD-186-0-1	12/28/2013 01:27	10	T1230590.D	RTX-1701 0.53(mm)
180-28384-25	SD-187-0-1	12/28/2013 01:58	10	T1230591.D	RTX-1701 0.53(mm)
180-28384-26	SD-188-0-1	12/28/2013 02:30	10	T1230592.D	RTX-1701 0.53(mm)
180-28384-27	SD-189-0-1	12/28/2013 03:01	10	T1230593.D	RTX-1701 0.53 (mm)
180-28384-28	SD-190-0-1	12/28/2013 03:33	10	T1230594.D	RTX-1701 0.53 (mm)
180-28384-29	SD-191-0-1	12/28/2013 04:04	10	T1230595.D	RTX-1701 0.53(mm)
180-28384-30	SD-192-0-1	12/28/2013 04:36	10.	T1230596.D	RTX-1701 0.53(mm)
180-28384-31	SD-193-0-1	12/28/2013 05:07	10	T1230597.D	RTX-1701 0.53(mm)
180-28384-32	SD-194-0-1	12/28/2013 05:39	10	T1230598.D	RTX-1701 0.53(mm)
MB 180-93333/1-C		12/28/2013 06:42	1	T1230600.D	RTX-1701 0.53(mm)
LCS 180-93333/2-C		12/28/2013 07:13	1	T1230601.D	RTX-1701 0.53(mm)
CCV 180-93552/58		12/28/2013 07:45	1	T1230602.D	RTX-1701 0.53(mm)
180-28384-13	SD-175-0-1	12/28/2013 08:16	50	T1230603.D	RTX-1701 0.53(mm)
180-28384-18	SD-180-0-1	12/28/2013 08:48	10	T1230604.D	RTX-1701 0.53(mm)
180-28384-21	SD-183-0-1	12/28/2013 10:23	2000	T1230607.D	RTX-1701 0.53(mm)
180-28384-19	SD-181-0-1	12/28/2013 11:58	20000	T1230610.D	RTX-1701 0.53 (mm)
180-28384-20	SD-182-0-3	12/28/2013 12:29	20000	T1230611.D	RTX-1701 0.53(mm)
CCV 180-93552/64		12/28/2013 15:07	1	T1230616.D	RTX-1701 0.53(mm)

### FORM VIII GC SEMI VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: ICRT 180-93552/10 Date Analyzed: 12/26/2013 22:40

Instrument ID: GC10 GC Column: RTX-1701 ID: 0.53(mm)

Lab File ID (Standard): T1230539.D Heated Purge: (Y/N) N

Calibration ID: 13174

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

				TCX	DCB
				RT #	RT #
INITIAL CALIBRATION S	SURROGATE	101-201-001-00-		5.66	17.60
UPPER LIMIT				5.71	17.65
LOWER LIMIT				5.61	17.55
LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	LAB FILE ID		
ICRT 180-93552/10		12/26/2013 22:40	T1230539.D	5.66	17.60
ICV 180-93552/20		12/27/2013 03:55	T1230549.D	0.00	0.00
180-28384-1	SD-163-0-1	12/27/2013 04:27	T1230550.D	5.66	17.59
180-28384-2	SD-164-0-1	12/27/2013 04:58	T1230551.D	5.66	17.59
180-28384-3	SD-165-0-1	12/27/2013 05:30	T1230552.D	5.66	17.59
180-28384-3 MS	SD-165-0-1 MS	12/27/2013 06:01	T1230553.D	5.66	17.59
180-28384-3 MSD	SD-165-0-1 MSD	12/27/2013 06:33	T1230554.D	5.66	17.60
180-28384-4	SD-166-0-1	12/27/2013 07:04	T1230555.D	5.66	17.60
180-28384-5	SD-167-0-1	12/27/2013 07:36	T1230556.D	5.66	17.59
180-28384-6	SD-168-0-1	12/27/2013 08:07	T1230557.D	5.66	17.57
180-28384-7	SD-169-0-1	12/27/2013 08:39	T1230558.D	5.66	17.59
180-28384-8	SD-170-0-1	12/27/2013 09:10	T1230559.D	5.66	17.59
CCV 180-93552/31		12/27/2013 13:22	T1230567.D	5.66	17.60
180-28384-12	SD-174-0-1	12/27/2013 14:26	T1230569.D	5.66	17.60
180-28384-14	SD-176-0-3	12/27/2013 15:28	T1230571.D	5.66	17.60
180-28384-15	SD-177-0-2	12/27/2013 16:00	T1230572.D	5.66	17.60
180-28384-16	SD-178-0-1	12/27/2013 16:31	T1230573.D	5.66	17.59
180-28384-17	SD-179-0-1	12/27/2013 17:03	T1230574.D	5.66	17.60
180-28384-9	SD-171-0-1	12/27/2013 19:09	T1230578.D	0.00	0.00
180-28384-10	SD-172-0-3	12/27/2013 19:40	T1230579.D	0.00	0.00
180-28384-11	SD-173-0-3	12/27/2013 20:12	T1230580.D	0.00	0.00
MB 180-93257/1-C		12/27/2013 21:15	T1230582.D	5.66	17.59
LCS 180-93257/2-C		12/27/2013 21:46	T1230583.D	5.66	17.59
CCV 180-93552/42		12/27/2013 22:18	T1230584.D	5.66	17.59
180-28384-22	SD-184-0-2	12/27/2013 23:21	T1230586.D	5.66	17.59
180-28384-23	SD-185-0-1	12/27/2013 23:52	T1230587.D	5.66	17.59
180-28384-23 MS	SD-185-0-1 MS	12/28/2013 00:24	T1230588.D	5.66	17.59

TCX = Tetrachloro-m-xylene

DCB = DCB Decachlorobiphenyl (Surr)

TCX RT Limit =  $\pm$  0.05 minutes of surrogate RT DCB RT Limit =  $\pm$  0.05 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8082A

### FORM VIII GC SEMI VOA ANALYTICAL SEQUENCE

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Sample No.: ICRT 180-93552/10 Date Analyzed: 12/26/2013 22:40

Instrument ID: GC10 GC Column: RTX-1701 ID: 0.53(mm)

Lab File ID (Standard): T1230539.D Heated Purge: (Y/N) N

Calibration ID: 13174

THE ANALYTICAL SEQUENCE OF BLANKS, SAMPLES, STANDARDS, MS/MSDs AND LCSs IS GIVEN BELOW:

				TCX	DCB
				RT #	RT #
INITIAL CALIBRATION S	BURROGATE		·	5.66	17.60
UPPER LIMIT				5.71	17.65
LOWER LIMIT	***************************************			5.61	17.55
LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	LAB FILE ID		
180-28384-23 MSD	SD-185-0-1 MSD	12/28/2013 00:55	T1230589.D	5.66	17.59
180-28384-24	SD-186-0-1	12/28/2013 01:27	T1230590.D	5.66	17.59
180-28384-25	SD-187-0-1	12/28/2013 01:58	T1230591.D	5.66	17.58
180-28384-26	SD-188-0-1	12/28/2013 02:30	T1230592.D	5.66	17.59
180-28384-27	SD-189-0-1	12/28/2013 03:01	T1230593.D	5.66	17.57
180-28384-28	SD-190-0-1	12/28/2013 03:33	T1230594.D	5.66	17.59
180-28384-29	SD-191-0-1	12/28/2013 04:04	T1230595.D	5.66	17.59
180-28384-30	SD-192-0-1	12/28/2013 04:36	T1230596.D	5.66	17.59
180-28384-31	SD-193-0-1	12/28/2013 05:07	T1230597.D	5.66	17.59
180-28384-32	SD-194-0-1	12/28/2013 05:39	T1230598.D	5.66	17.59
MB 180-93333/1-C		12/28/2013 06:42	T1230600.D	5.66	17.59
LCS 180-93333/2-C		12/28/2013 07:13	T1230601.D	5.66	17.59
CCV 180-93552/58		12/28/2013 07:45	T1230602.D	5.66	17.59
180-28384-13	SD-175-0-1	12/28/2013 08:16	T1230603.D	5.66	17.59
180-28384-18	SD-180-0-1	12/28/2013 08:48	T1230604.D	5.66	17.59
180-28384-21	SD-183-0-1	12/28/2013 10:23	T1230607.D	0.00	0.00
180-28384-19	SD-181-0-1	12/28/2013 11:58	T1230610.D	0.00	0.00
180-28384-20	SD-182-0-3	12/28/2013 12:29	T1230611.D	0.00	0.00
CCV 180-93552/64		12/28/2013 15:07	T1230616.D	5.66	17.60

TCX = Tetrachloro-m-xylene

DCB = DCB Decachlorobiphenyl (Surr)

TCX RT Limit =  $\pm$  0.05 minutes of surrogate RT DCB RT Limit =  $\pm$  0.05 minutes of surrogate RT

# Column used to flag values outside QC limits

FORM VIII 8082A

### GC SEMI VOA INITIAL CALIBRATION DATA EXTERNAL STANDARD CURVE EVALUATION FORM VI

Analy Batch No.: 93552 Calibration ID: 13168 Heated Purge: (Y/N) 12/26/2013 17:57 ID: 0.53(mm) Calibration End Date: Job No.: 180-28384-1 GC Column: RTX-1701 12/26/2013 17:57 Lab Name: TestAmerica Pittsburgh Calibration Start Date: Instrument ID: GC10 SDG No.:

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
L ( C ( )	100 005 07	i control

	OX COD								
_	*KSU OK COU	0.0	0.0	20.0	0.0	0.0	0.0	0.0	20.02
%RSD # M	Y,	20	. 2(	2(	2(	2(	5(	75	20
# MIN CF									
	M2								
COEFFICIENT	M1	186256.000	115178.000	358064.000	267766.000	421038.000	306508.000	582504.000	362774.000
CURVE	TYPE	Ave							
		_							
CF			-						
	LVL 1	186256	115178	358064	267766	421038	306508	582504	362774
ANALYTE		PCB-1221 Peak 1	PCB-1221 Peak 2	PCB-1221 Peak 3	PCB-1254 Peak 1	PCB-1254 Peak 2	PCB-1254 Peak 3	PCB-1254 Peak 4	PCB-1254 Peak 5

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### GC SEMI VOA INITIAL CALIBRATION DATA EXTERNAL STANDARD CURVE EVALUATION FORM VI

93552 Z Heated Purge: (Y/N) Analy Batch No.: ID: 0.53(mm) Job No.: 180-28384-1 GC Column: RTX-1701 Lab Name: TestAmerica Pittsburgh Instrument ID: GC10 SDG No.:

Calibration ID: 13169

12/26/2013 18:28

Calibration End Date:

12/26/2013 18:28

Calibration Start Date:

LAB FILE ID: T1230531.D LAB SAMPLE ID: IC 180-93552/2 Calibration Files: LEVEL: Level 1

MIN R^2	OK COD					
R^2 #	*KSD OK COD	-				
MAX	*KSD	20.0	20.0	20.0	20.0	20.0
%RSD #						
MIN CF   %RSD						
#		_				
	M2					
COEFFICIENT	M1	103436.000	78870.0000	106964.000	151424.000	123748.000
	м					
CURVE	TYPE	Ave	Ave	Ave	Ave	Ave
CF						
	LVL 1		78870	106964	151424	123748
ANALYTE		PCB-1232 Peak 1	PCB-1232 Peak 2	PCB-1232 Peak 3	PCB-1232 Peak 4	PCB-1232 Peak 5

# GC SEMI VOA INITIAL CALIBRATION DATA EXTERNAL STANDARD CURVE EVALUATION

Lab Name: Tes	Lab Name: TestAmerica Pittsburgh	Job No.: 180-28384-1	180-28384-1		Analy Batch No.:	93552	
SDG No.:						-	
Instrument ID: GC10	): GC10	GC Column:	RTX-1701	ID: 0.53 (mm)	Heated Purge: (Y/N)	Z	

Calibration End Date: 12/26/2013 19:00

12/26/2013 19:00

Calibration ID: 13170

Calibration Files:

Calibration Start Date:

LEVEL: LAB SAMPLE ID: LAB FILE ID: Level 1 IC 180-93552/3 T1230532.D

ANALYTE	CF	CURVE	COEFFICIENT		MIN CF	MIN CF %RSD #	MAX	R^2	#
1 11 21 21 21 22 22 22 22 22 22 22 22 22	LVL 1	TYPE	B M1 M2	- 2			%RSD	OR COD	OR COD
PCB-1242 Peak 1	445914	Ave	445914.000				20.0		
2 Peak 2	181944	Ave	181944.000				20.0		
2 Peak 3	207238	Ave	207238.000				20.0		
PCB-1242 Peak 4	262994	Ave	262994.000				20.0		
-1242 Peak 5	150326	Ave	150326.000				20.0		

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### GC SEMI VOA INITIAL CALIBRATION DATA EXTERNAL STANDARD CURVE EVALUATION FORM VI

Job No.: 180-28384-1 Lab Name: TestAmerica Pittsburgh

Instrument ID: GC10

12/26/2013 19:31

Calibration Start Date:

SDG No.:

Calibration ID: 13171 Heated Purge: (Y/N) 12/26/2013 19:31 ID: 0.53(mm) Calibration End Date: GC Column: RTX-1701

Analy Batch No.: 93552

Calibration Files:

LAB FILE ID: T1230533.D LAB SAMPLE ID: IC 180-93552/4 LEVEL: Level 1

MAX R^2 # MIN R^2		20.0	20.0	20.0	20.0	20.0
%RSD #						
# MIN CF		-				
	M2					
COEFFICIENT	M1	367610.000	138678.000	416140.000	327236.000	196130,000
SURVE	TYPE					
O.F.	LVL 1			416140 Ave	327236 A	196130 Ave
ANALYTE		PCB-1248 Peak 1	CB-1248 Peak 2	PCB-1248 Peak 3	CB-1248 Peak 4	PCB-1248 Peak 5

### GC SEMI VOA INITIAL CALIBRATION DATA EXTERNAL STANDARD CURVE EVALUATION FORM VI

93552 Calibration ID: 13174 Heated Purge: (Y/N) Analy Batch No.: 12/27/2013 00:15 ID: 0.53(mm) Calibration End Date: Job No.: 180-28384-1 GC Column: RTX-1701 12/26/2013 21:06 Lab Name: TestAmerica Pittsburgh Calibration Start Date: Instrument ID: GC10 SDG No.:

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 180-93552/7	T1230536.D
Level 2	IC 180-93552/8	T1230537.D
Level 3	IC 180-93552/9	T1230538.D
Level 4	ICRT 180-93552/10	T1230539.D
Level 5	IC 180-93552/11	T1230540.D
Level 6	IC 180-93552/12	T1230541.D
Level 7	IC 180-93552/13	T1230542.D

ANALYTE		CE		<u>ੂ</u>	CURVE	COEFFICIENT		# MIN CF	%RSD #	MAX	MIM #
;	LVL 1 LVL 5	LVL 2 LVL 6	LVL 3 LVL 7	LVL 4	e HAAA.	M1	M2			*Kad OK COD	Z Z Z
PCB-1016 Peak 1	316100	321920	324650	272626 Ave	Ф	298692.607			7.3	20.0	
PCB-1016 Peak 2	721900	669580		557236 Ave	ō.	616141.679			10.6	20.0	
PCB-1016 Peak 3	922400	855600	887920	757466 Ave	Ф	854659.179			6.3	20.0	
PCB-1016 Peak 4	383100	382100	382720	315518 Ave	, a	357419.714			7.4	20.0	
PCB-1016 Peak 5	301000	291900	292005	239116 Ave	ø	275455.714		-	8.1	20.0	<u> </u>
PCB-1260 Peak 1	551100	538120	525755 440183	449752 Ave	φ.	489514.786			9.6	20.0	
PCB-1260 Peak 2	615200	624580	612015	518906 Ave	ē.	566016.857			8.0	20.0	
PCB-1260 Peak 3	601200	595100		491290 Ave	ė	540242.786			1.6	20.0	-
PCB-1260 Peak 4	431700	441180	434045	368872 Ave	ė.	398970.179			8	20.0	<u> </u>
PCB-1260 Peak 5	896200	879360		785820 Ave	p.	849164.964			5.1	20.0	<u>:</u> 
Tetrachloro-m-xylene	28846000	23855600	24641400	21955760 Ave	ų.	24259487.9			9.1	20.0	
DCB Decachlorobiphenyl (Surr)	9714000	8719600 8018880	9245100 7675175	8010200 Ave	Đ	8526847.86		-	8	20.0	

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: ICV 180-93552/14 Calibration Date: 12/27/2013 00:46

Instrument ID: GC10 Calib Start Date: 12/26/2013 17:57

GC Column: RTX-1701 ID: 0.53(mm) Calib End Date: 12/26/2013 17:57

Lab File ID: T1230543.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE CF	CF	MIN CF	CALC AMOUNT	SPIKE AMOUNT	용D	MAX %D
PCB-1221 Peak 1	Ave	186256	180814		0.485	0.500	-2.9	20.0
PCB-1221 Peak 2	Ave	115178	113616		0.493	0.500	-1.4	20.0
PCB-1221 Peak 3	Ave	358064	354118		0.494	0.500	-1.1	20.0
PCB-1254 Peak 1	Ave	267766	290136		0.542	0.500	8.4	20.0
PCB-1254 Peak 2	Ave	421038	448990		0.533	0.500	6.6	20.0
PCB-1254 Peak 3	Ave	306508	339866		0.554	0.500	10.9	20.0
PCB-1254 Peak 4	Ave	582504	648740		0.557	0.500	11.4	20.0
PCB-1254 Peak 5	Ave	362774	388648		0.536	0.500	7.1	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: <u>ICV 180-93552/15</u> Calibration Date: 12/27/2013 01:18

Instrument ID: GC10 Calib Start Date: 12/26/2013 18:28

GC Column: RTX-1701 ID: 0.53(mm) Calib End Date: 12/26/2013 18:28

Lab File ID: T1230544.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE CF	CF	MIN CF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
PCB-1232 Peak 1	Ave	103436	105430		0.510	0.500	1.9	20.0
PCB-1232 Peak 2	Ave	78870	80908		0.513	0.500	2.6	20.0
PCB-1232 Peak 3	Ave	106964	109506		0.512	0.500	2.4	20.0
PCB-1232 Peak 4	Ave	151424	153464		0.507	0.500	1.3	20.0
PCB-1232 Peak 5	Ave	123748	127758		0.516	0.500	3.2	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: ICV 180-93552/16 Calibration Date: 12/27/2013 01:49

Instrument ID: GC10 Calib Start Date: 12/26/2013 19:00

GC Column: RTX-1701 ID: 0.53(mm) Calib End Date: 12/26/2013 19:00

Lab File ID: T1230545.D Conc. Units: ng/uL

	ANALYTE	CURVÉ TYPE	AVE CF	CF	MIN CF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
PCB-1242	Peak 1	Ave	445914	473648		0.531	0.500	6.2	20.0
PCB-1242	Peak 2	Ave	181944	180526		0.496	0.500	-0.8	20.0
PCB-1242	Peak 3	Ave	207238	204500		0.493	0.500	-1.3	20.0
PCB-1242	Peak 4	Ave	262994	268914		0.511	0.500	2.3	20.0
PCB-1242	Peak 5	Ave	150326	146902		0.489	0.500	-2.3	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: ICV 180-93552/17 Calibration Date: 12/27/2013 02:21

Instrument ID: GC10 Calib Start Date: 12/26/2013 19:31

GC Column: RTX-1701 ID: 0.53(mm) Calib End Date: 12/26/2013 19:31

Lab File ID: T1230546.D Conc. Units: ng/uL

	ANALYTE	CURVE TYPE	AVE CF	CF	MIN CF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
PCB-1248 E	Peak 1	Ave	367610	304126		0.414	0.500	-17.3	20.0
PCB-1248 E	Peak 2	Ave	138678	116780		0.421	0.500	-15.8	20.0
PCB-1248 E	Peak 3	Ave	416140	354832		0.426	0.500	-14.7	20.0
PCB-1248 E	Peak 4	Ave	327236	297900		0.455	0.500	-9.0	20.0
PCB-1248 E	Peak 5	Ave	196130	182354		0.465	0.500	-7.0	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: ICV 180-93552/20 Calibration Date: 12/27/2013 03:55

Instrument ID: GC10 Calib Start Date: 12/26/2013 21:06

GC Column: RTX-1701 ID: 0.53(mm) Calib End Date: 12/27/2013 00:15

Lab File ID: T1230549.D Conc. Units: ng/uL

ANALYTE	CURVE	AVE CF	CF	MIN CF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
PCB-1016 Peak 1	Ave	298693	253758		0.425	0.500	-15.0	20.0
PCB-1016 Peak 2	Ave	616142	546314		0.443	0.500	-11.3	20.0
PCB-1016 Peak 3	Ave	854659	748164		0.438	0.500	-12.5	20.0
PCB-1016 Peak 4	Ave	357420	314542		0.440	0.500	-12.0	20.0
PCB-1016 Peak 5	Ave	275456	240406		0.436	0.500	-12.7	20.0
PCB-1260 Peak 1	Ave	489515	469866		0.480	0.500	-4.0	20.0
PCB-1260 Peak 2	Ave	566017	537872.		0.475	0.500	-5.0	20.0
PCB-1260 Peak 3	Ave	540243	507318		0.470	0.500	-6.1	20.0
PCB-1260 Peak 4	Ave	398970	388902		0.487	0.500	-2.5	20.0
PCB-1260 Peak 5	Ave	849165	816856		0.481	0.500	-3.8	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCV 180-93552/31 Calibration Date: 12/27/2013 13:22

Instrument ID: GC10 Calib Start Date: 12/26/2013 21:06

GC Column: RTX-1701 ID: 0.53(mm) Calib End Date: 12/27/2013 00:15

Lab File ID: T1230567.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE CF	CF	MIN CF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
PCB-1016 Peak 1	Ave	298693	309384		1.04	1.00	3.6	20.0
PCB-1016 Peak 2	Ave	616142	623214		1.01	1.00	1.1	20.0
PCB-1016 Peak 3	Ave	854659	889676		1.04	1.00	4.1	20.0
PCB-1016 Peak 4	Ave	357420	373856		1.05	1.00	4.6	20.0
PCB-1016 Peak 5	Ave	275456	288706		1.05	1.00	4.8	20.0
PCB-1260 Peak 1	Ave	489515	481278		0.983	1.00	-1.7	20.0
PCB-1260 Peak 2	Ave	566017	557420		0.985	1.00	-1.5	20.0
PCB-1260 Peak 3	Ave	540243	525912		0.973	1.00	-2.7	20.0
PCB-1260 Peak 4	Ave	398970	385777		0.967	1.00	-3.3	20.0
PCB-1260 Peak 5	Ave	849165	824971		0.972	1.00	-2.8	2.0.0
Tetrachloro-m-xylene	Ave	24259488	25277800		0.0521	0.0500	4.2	20.0
DCB Decachlorobiphenyl (Surr)	Ave	8526848	7584760		0.0445	0.0500	-11.0	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCV 180-93552/42 Calibration Date: 12/27/2013 22:18

Instrument ID: GC10 Calib Start Date: 12/26/2013 21:06

GC Column: RTX-1701 ID: 0.53(mm) Calib End Date: 12/27/2013 00:15

Lab File ID: T1230584.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE CF	CF	MIN CF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
PCB-1016 Peak 1	Ave	298693	313303		1.05	1.00	4.9	20.0
PCB-1016 Peak 2	Ave	616142	623305		1.01	1.00	1.2	20.0
PCB-1016 Peak 3	Ave	854659	908124		1.06	1.00	6.3	20.0
PCB-1016 Peak 4	: Ave	357420	379352		1.06	1.00	6.1	20.0
PCB-1016 Peak 5	Ave	275456	293450		1.07	1.00	6.5	20.0
PCB-1260 Peak 1	Ave	489515	493291		1.01	1.00	0.8	20.0
PCB-1260 Peak 2	Ave	566017	577675		1.02	1.00	2.1	20.0
PCB-1260 Peak 3	Ave	540243	538644		0.997	1.00	-0.3	20.0
PCB-1260 Peak 4	Ave	398970	398658		0.999	1.00	-0.0	20.0
PCB-1260 Peak 5	Ave	849165	844400		0.994	1.00	-0.6	20.0
Tetrachloro-m-xylene	Ave	24259488	25280260		0.0521	0.0500	4.2	20.0
DCB Decachlorobiphenyl (Surr)	Ave	8526848	7416580		0.0435	0.0500	-13.0	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCV 180-93552/58 Calibration Date: 12/28/2013 07:45

Instrument ID: GC10 Calib Start Date: 12/26/2013 21:06

GC Column: RTX-1701 ID: 0.53(mm) Calib End Date: 12/27/2013 00:15

Lab File ID: T1230602.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE CF	CF	MIN CF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
PCB-1016 Peak 1	Ave	298693	292456		0.490	0.500	-2.1	20.0
PCB-1016 Peak 2	Ave	616142	601886		0.488	0.500	-2.3	20.0
PCB-1016 Peak 3	Ave	854659	829186		0.485	0.500	-3.0	20.0
PCB-1016 Peak 4	Ave	357420	352062		0.493	0.500	-1.5	20.0
PCB-1016 Peak 5	Ave	275456	272830		0.495	0.500	-1.0	20.0
PCB-1260 Peak 1	Ave	489515	477412		0.488	0.500	-2.5	20.0
PCB-1260 Peak 2	Ave	566017	547626		0.484	0.500	-3.2	20.0
PCB-1260 Peak 3	Ave	540243	51798.4		0.479	0.500	-4.1	20.0
PCB-1260 Peak 4	Ave	398970	382532		0.479	0.500	-4.1	20.0
PCB-1260 Peak 5	Ave	849165	775262		0.456	0.500	-8.7	20.0
Tetrachloro-m-xylene	Ave	24259488	23823440		0.0246	0.0250	-1.8	20.0
DCB Decachlorobiphenyl (Surr)	Ave	8526848	6601840		0.0194	0.0250	-22.6*	20.0

Lab Name: TestAmerica Pittsburgh Job No.: 180-28384-1

SDG No.:

Lab Sample ID: CCV 180-93552/64 Calibration Date: 12/28/2013 15:07

Instrument ID: GC10 Calib Start Date: 12/26/2013 21:06

GC Column: RTX-1701 ID: 0.53(mm) Calib End Date: 12/27/2013 00:15

Lab File ID: T1230616.D Conc. Units: ng/uL

ANALYTE	CURVE TYPE	AVE CF	CF	MIN CF	CALC AMOUNT	SPIKE AMOUNT	% & D	MAX %D
PCB-1016 Peak 1	Ave	298693	281292		0.471	0.500	-5.8	20.0
PCB-1016 Peak 2	Ave	616142	. 581766		0.472	0.500	-5.6	20.0
PCB-1016 Peak 3	Ave	854659	800442		0.468	0.500	-6.3	20.0
PCB-1016 Peak 4	Ave	357420	331934		0.464	0.500	-7.1	20.0
PCB-1016 Peak 5	Ave	275456	256490		0.466	0.500	-6.9	20.0
PCB-1260 Peak 1	Ave	489515	462622		0.473	0.500	-5.5	20.0
PCB-1260 Peak 2	Ave	566017	531850		0.470	0.500	-6.0	20.0
PCB-1260 Peak 3	Ave	540243	498412		0.461	0.500	-7.7	20.0
PCB-1260 Peak 4	Ave	398970	375246		0.470	0.500	-5.9	20.0
PCB-1260 Peak 5	Ave	849165	761052		0.448	0.500	-10.4	20.0
Tetrachloro-m-xylene	Ave	24259488	22764120		0.0235	0.0250	-6.2	20.0
DCB Decachlorobiphenyl (Surr)	Ave	8526848	6472600		0.0190	0.0250	-24.1*	20.0

Data File: T1230550.D

Date: 27-DEC-2013 04:27

Client ID:

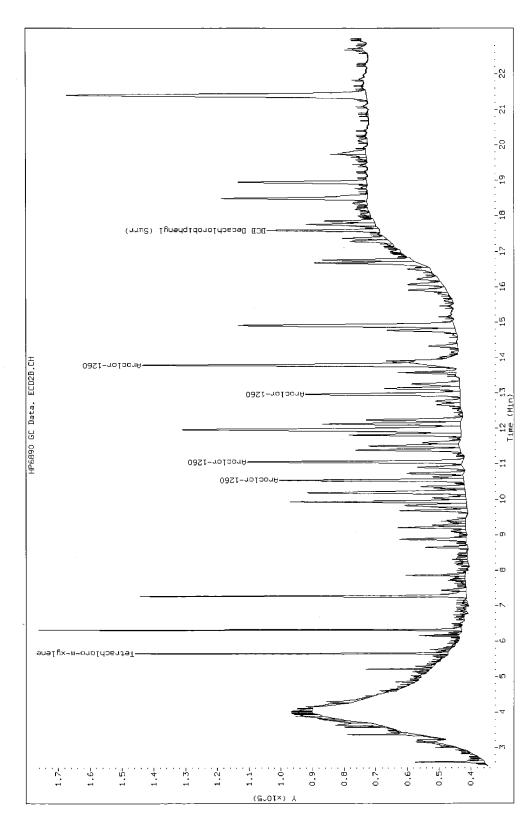
Sample Info: 12273B8082ALL.b

Instrument: gc10.i

SD-163-0-

Operator: 402360





Data File: T1230551.D

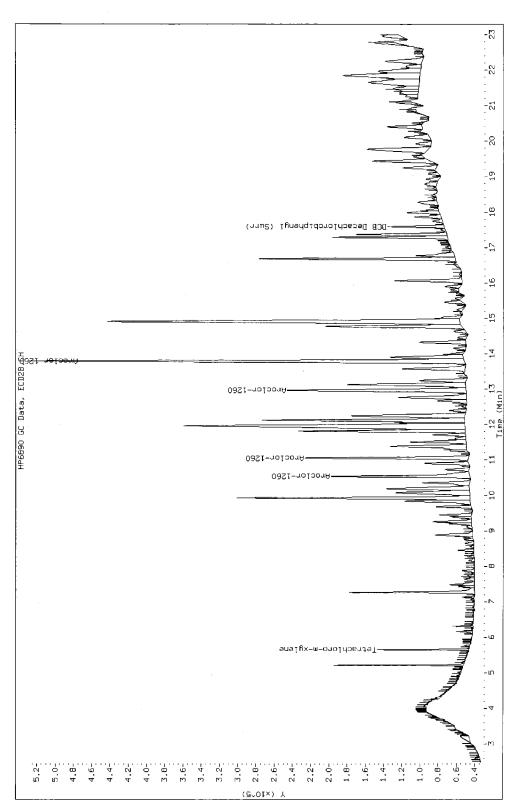
1-0-191-as Date: 27-DEC-2013 04:58

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i

Operator: 402360



Data File: T1230552.D

Date: 27-DEC-2013 05:30

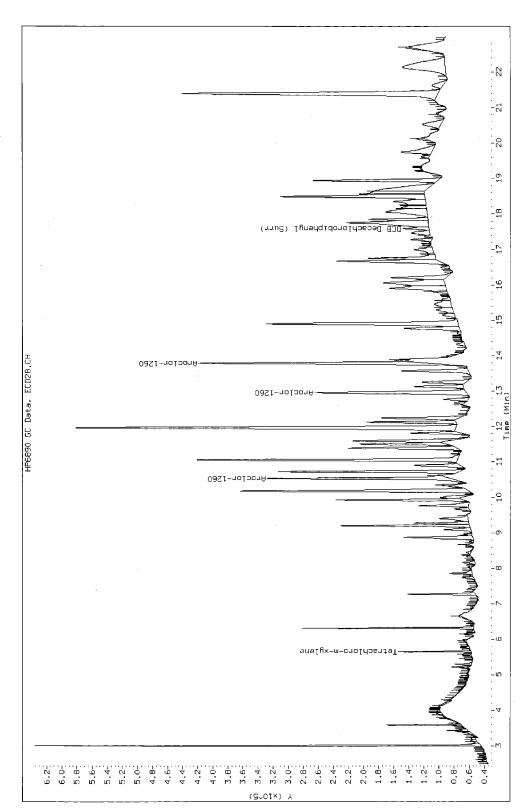
SD-165-0-

Client ID:

Instrument: gc10.i

Sample Info: 12273B8082ALL.b

Operator: 402360



Data File: T1230555.D

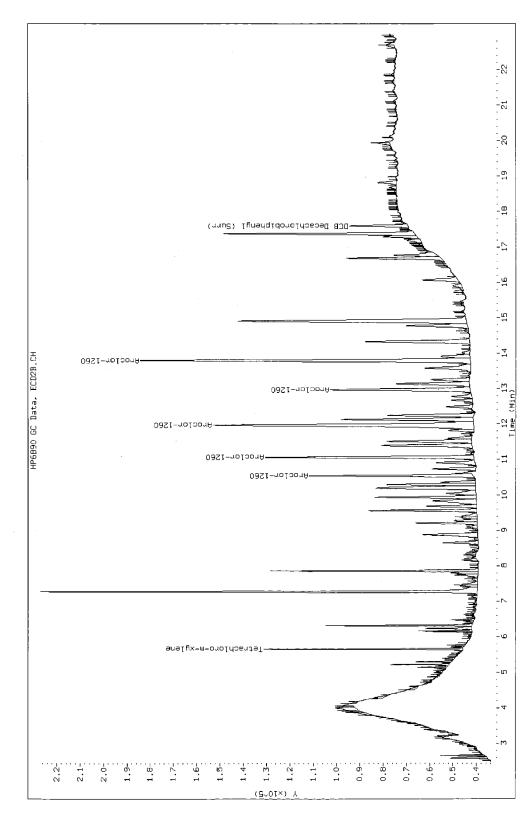
Date: 27-DEC-2013 07:04

Client ID:

Sample Info: 12273B8082ALL.b

SD-166-0-

Instrument: gc10.i Operator: 402360



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Data File: T1230556.D

Date: 27-DEC-2013 07:36

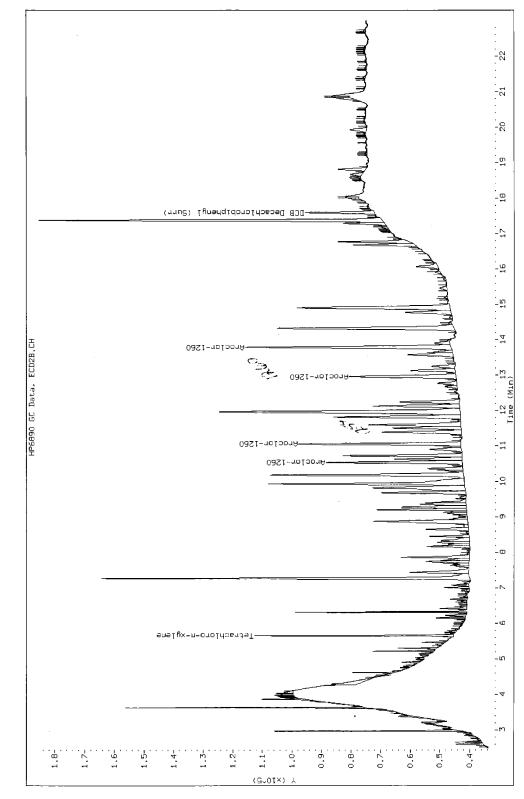
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Sample Info: 12273B8082ALL.b

SD-167-0-1

Operator: 402360

Instrument: gc10.i



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Data File: T1230557.D

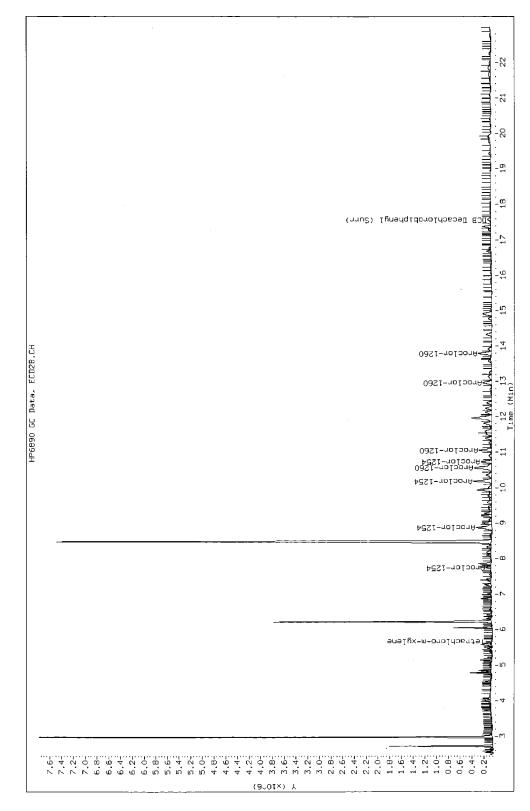
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1-0-891-as

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Sample Info: 12273B8082ALL.b

Instrument: gc10.i



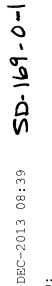
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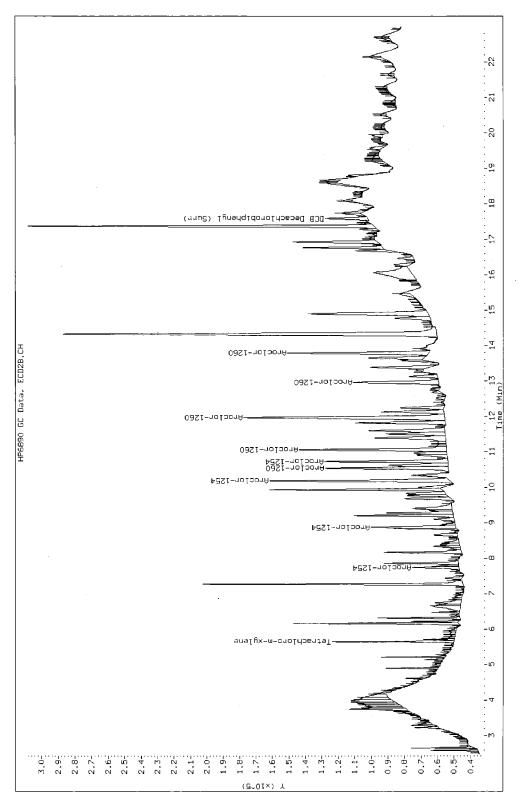
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Instrument: gc10.i





Data File: T1230559.D

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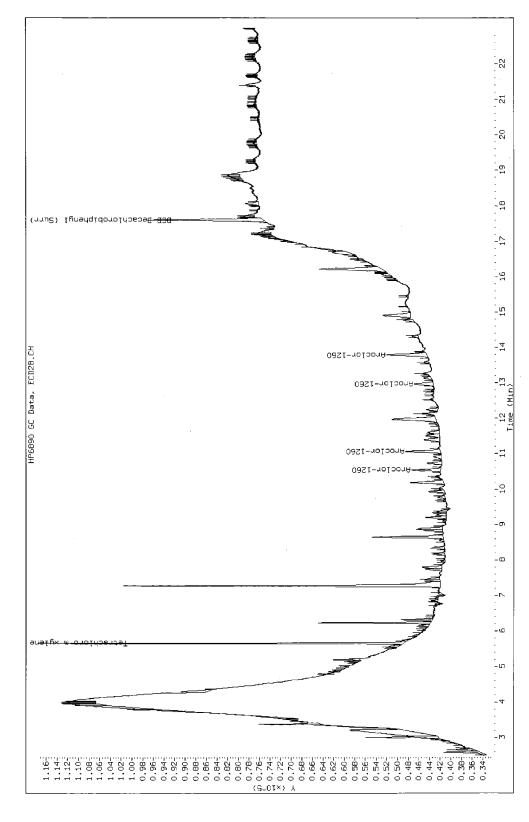
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1-0-0L1-QS

Instrument: gc10.i

Sample Info: 12273B8082ALL.b



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Data File: T1230578.D

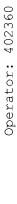
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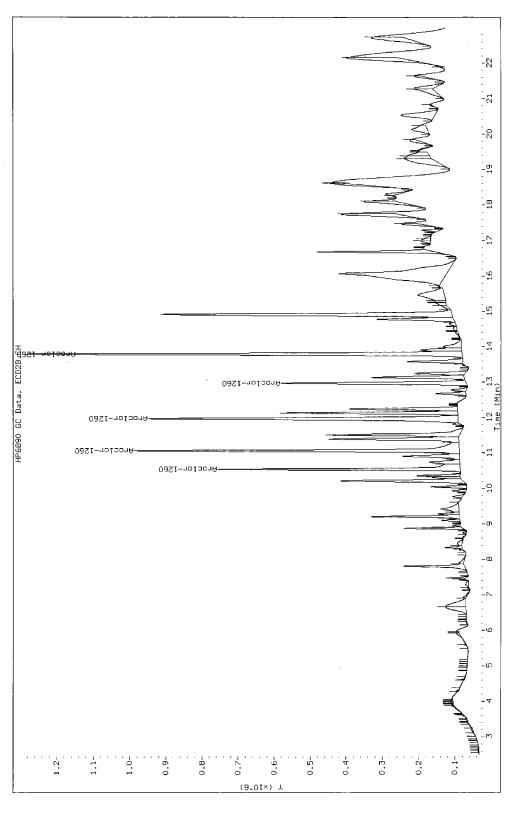
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Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i





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Data File: T1230579.D

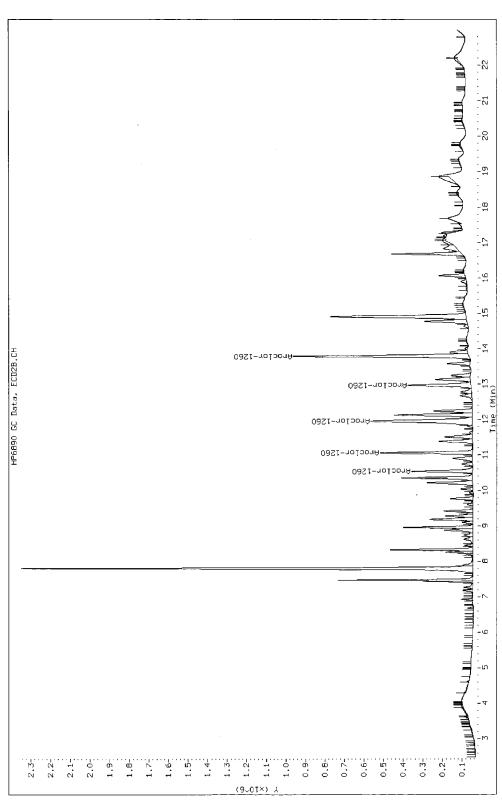
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SD-172-0-3

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Instrument: gc10.i

Sample Info: 12273B8082ALL.b



Data File: T1230580.D

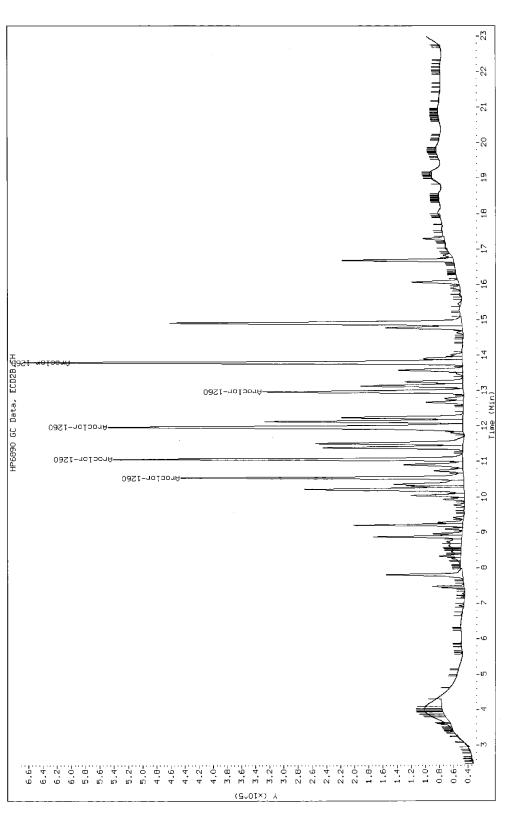
Date: 27-DEC-2013 20:12

sp-173-0-3

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



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Data File: T1230569.D

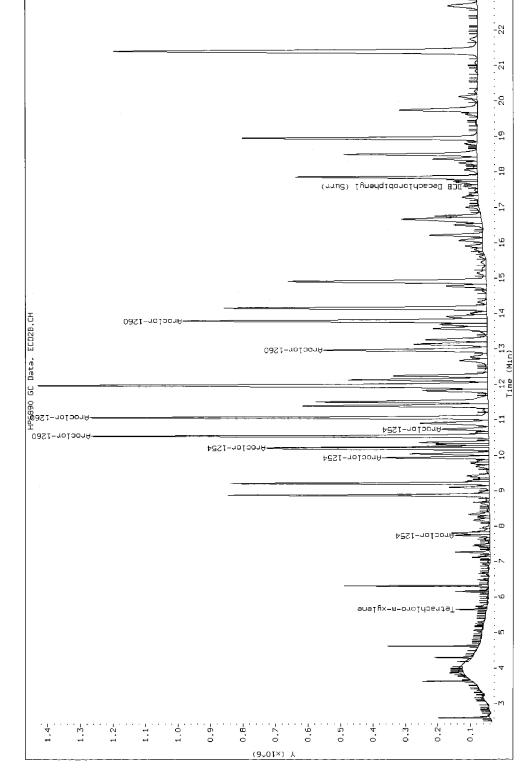
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1-0- HL1 -05

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



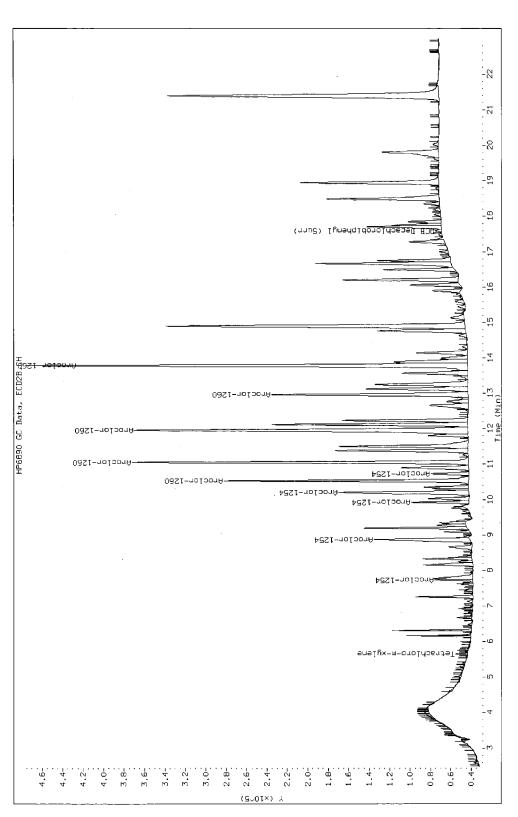
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Date: 28-DEC-2013 08:16 SD-175-0-1

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



Data File: T1230571.D

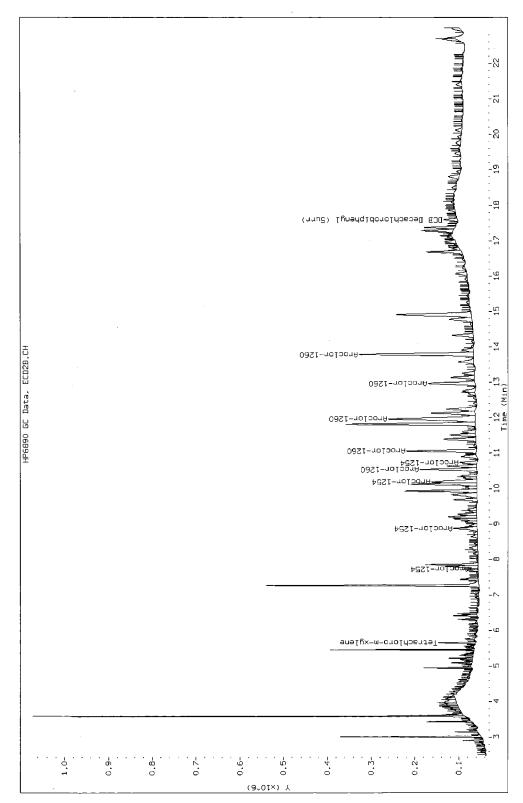
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Client ID:

Instrument: gc10.i

Sample Info: 12273B8082ALL.b



Data File: T1230572.D

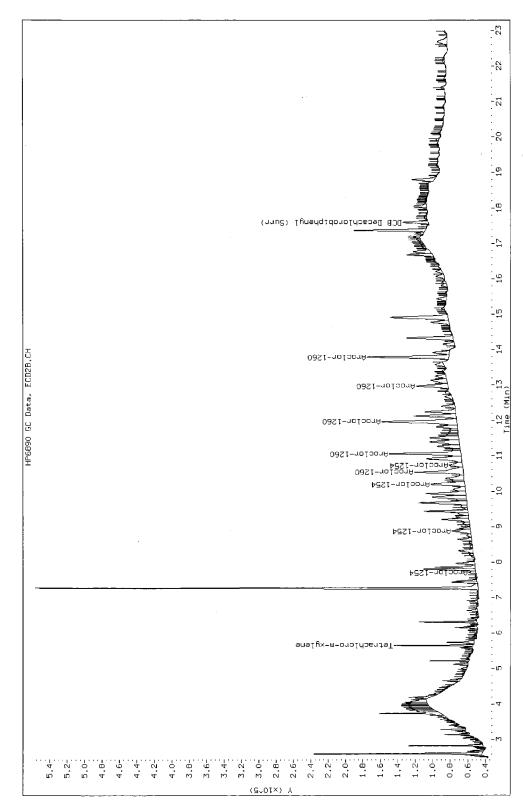
c-0-LL1-as

Date: 27-DEC-2013 16:00

Client ID:

Instrument: gc10.i

Sample Info: 12273B8082ALL.b



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Data File: T1230573.D

Date: 27-DEC-2013 16:31

Client ID:

SD-178-0

Instrument: gc10.i Operator: 402360

Sample Info: 12273B8082ALL.b

17 18 19 20 21 22 OCB Decachlorobiphenyl (Surv) -19 -記 9 10 11 12 13 14 Time (Min) HP6890 GC Data, ECD2B,CH 09S1-roloorA 06S1-701aoń£ 08S1-701⊃07A 401007A-03S1-701307A Arocloon-1254 5 0.81-10 0.78-X 0.75-0.42-0,93-0.90-0.84-0.69-0.66 0.63-0,60 0.54-0.48-1.14- $1.11^{-1}$ 1.08-1.05-1.02-0.99--96-0 0.72-0.57-0.51-0,45-1.17-

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Data File: T1230574.D

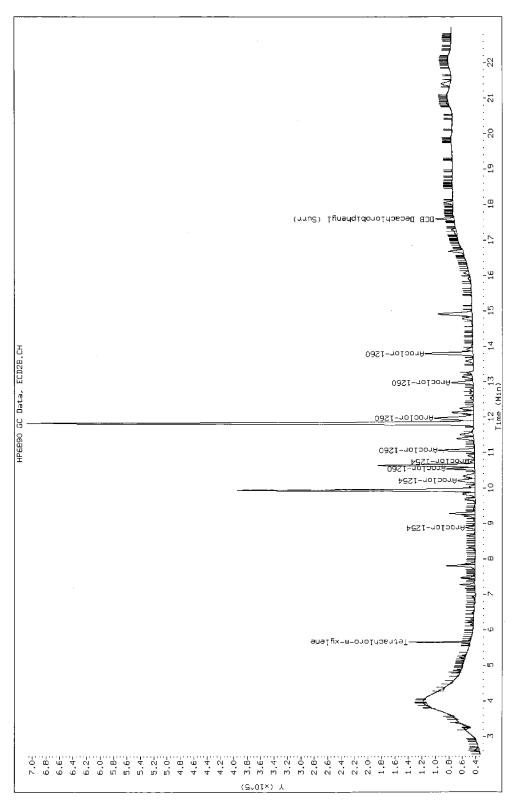
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1-0-6L1-0S

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Instrument: gc10.i

Sample Info: 12273B8082ALL.b



Data File: T1230604.D

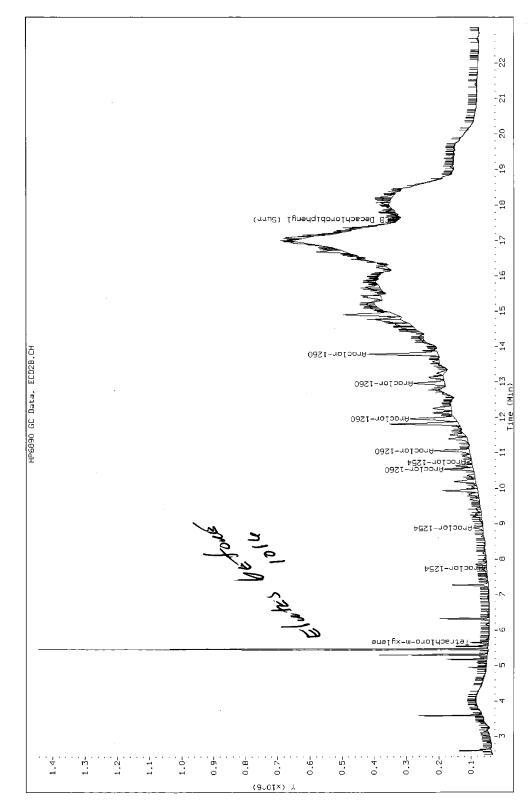
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-0-081-ds

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



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Data File: T1230610.D

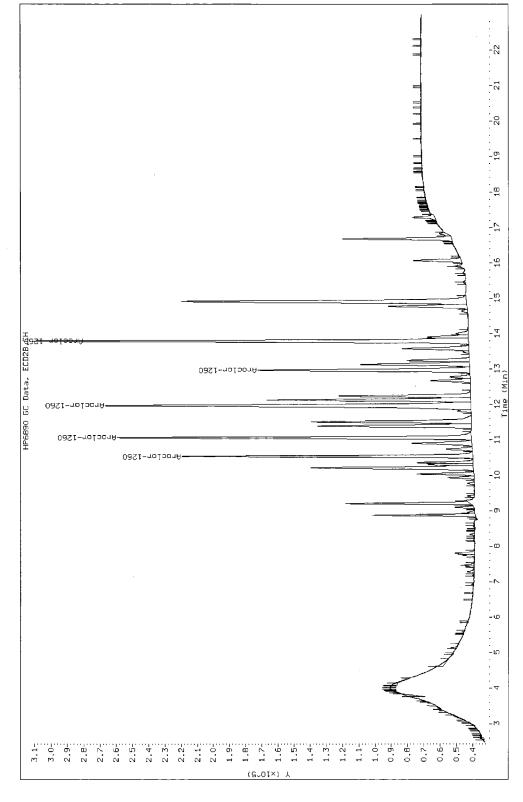
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Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i

1-0-181-OS



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Data File: T1230611.D

240. 28-PEC-2013 12.00

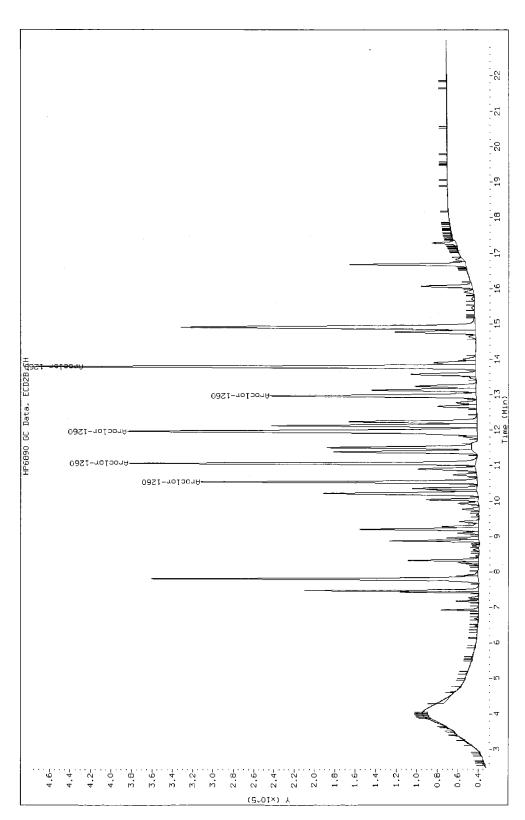
SD- 189-0-3

Date: 28-DEC-2013 12:29

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



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Data File: T1230607.D

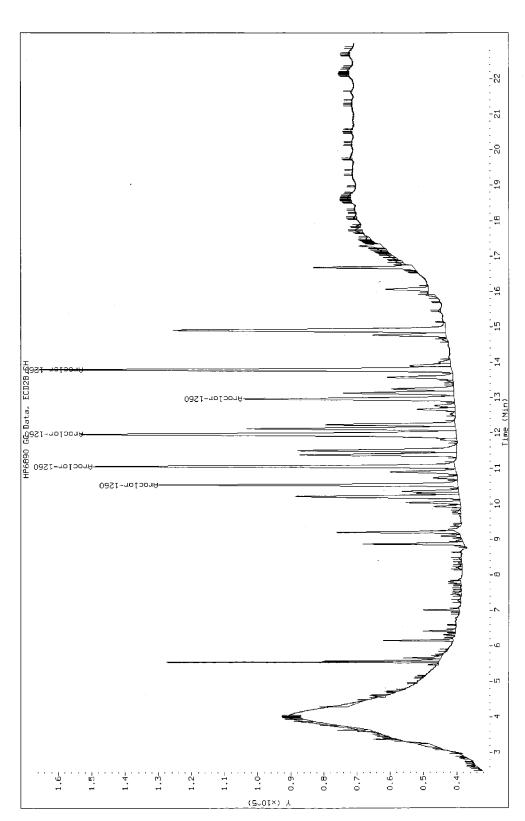
Date: 28-DEC-2013 10:23

SD-183-0-1

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



Data File: T1230586.D

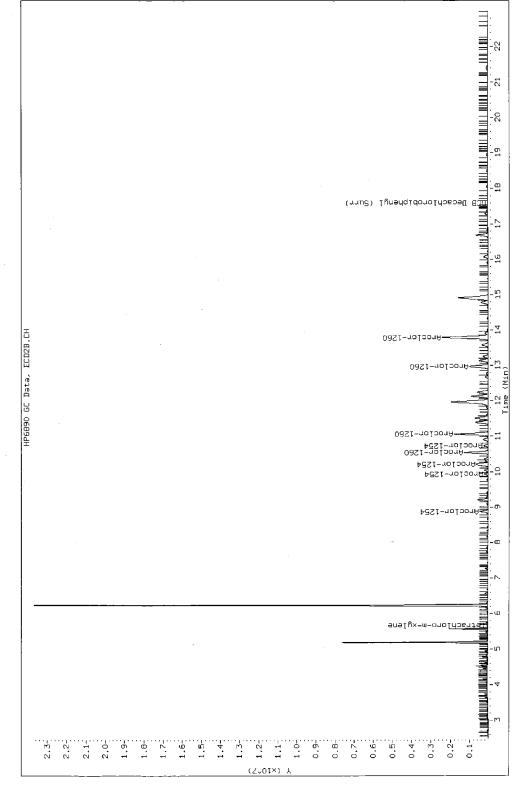
SD-184-0-2

Date: 27-DEC-2013 23:21

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



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Data File: T1230587.D

Date: 27-DEC-2013 23:52

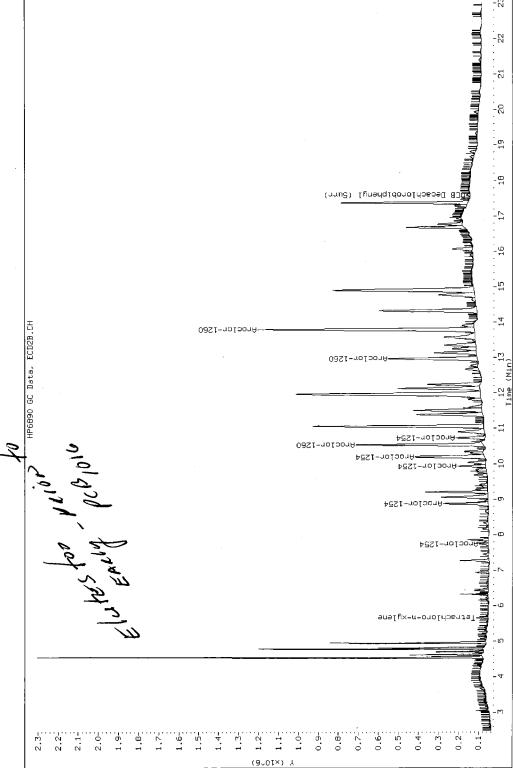
SD-185-0-1

Client ID:

Instrument: gc10.i

Operator: 402360

Sample Info: 12273B8082ALL.b



Data File: T1230590.D

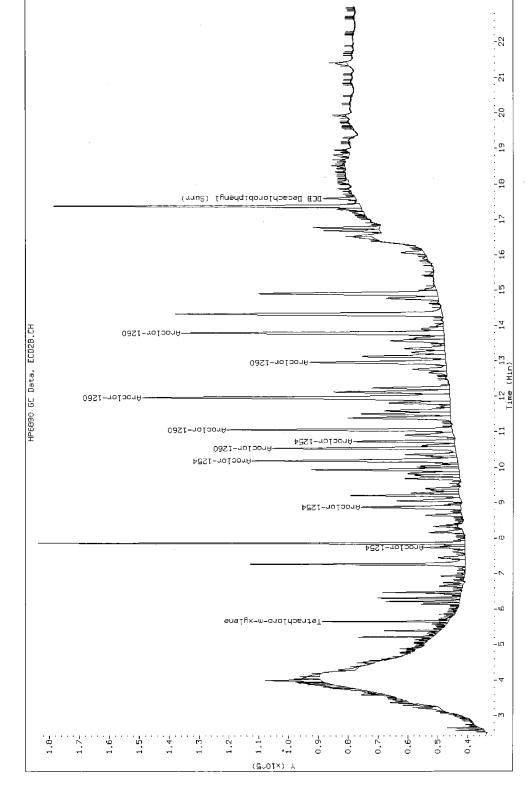
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50-186-0-1

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



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Data File: T1230591.D

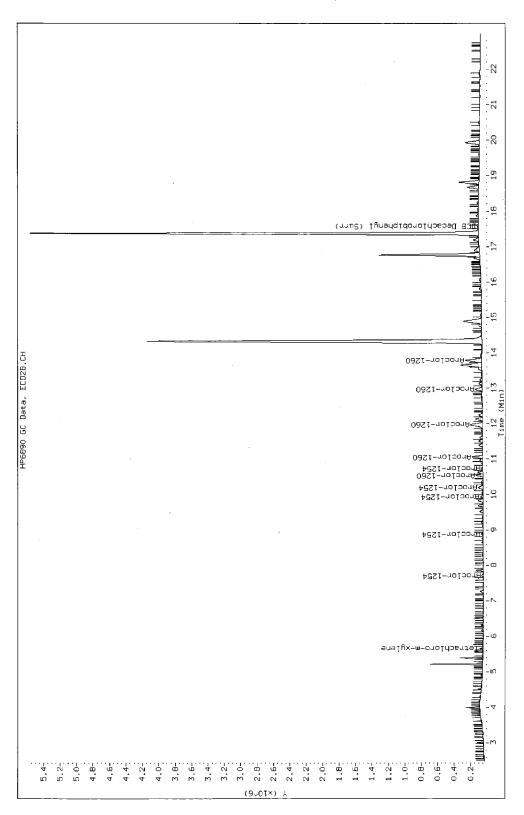
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SD-181-02

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



Data File: T1230592.D

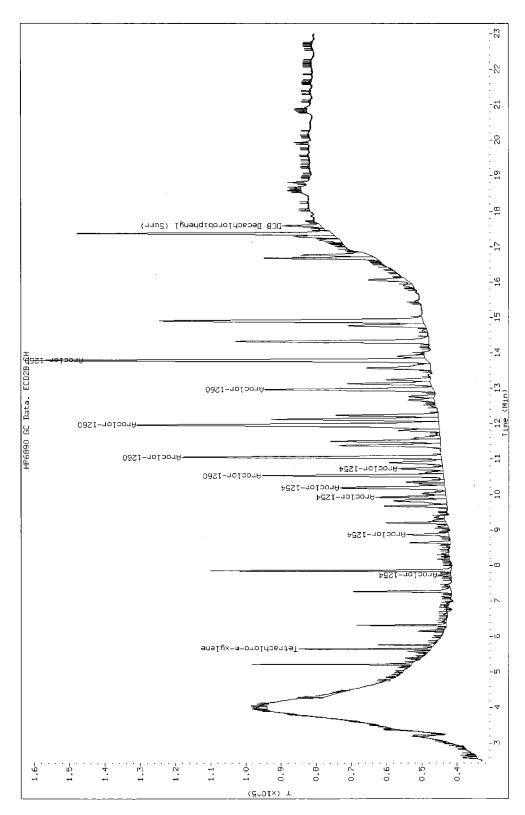
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1-2-881-05

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Sample Info: 12273B8082ALL.b

Instrument: gc10.i



Data File: T1230593.D

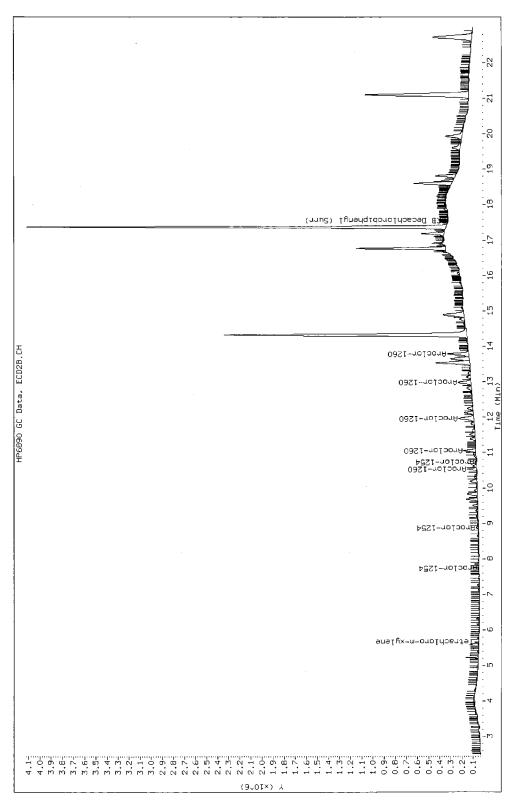
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1-0-1891-05

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



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Data File: T1230594.D

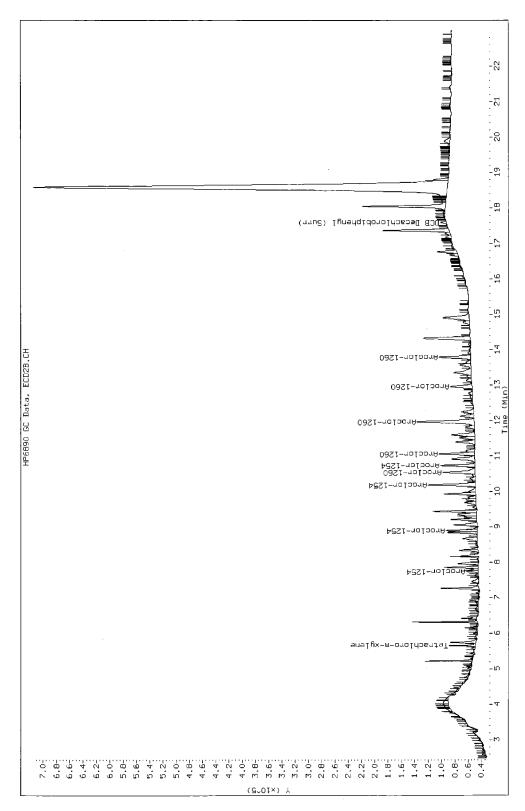
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1-0-061-05

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



Data File: T1230596.D

Date: 28-DEC-2013 04:36

Client ID:

SD-192-0-

Instrument: gc10.i

Operator: 402360

Sample Info: 12273B8082ALL.b

. 25 . . 21 19 20 -1 (vru2) Iynahqidoroidosəd<del>-də</del> . 17 . 16 -11 -12 -<u>4</u> HP6890 GC Data, ECD2B.CH 08S1-roloorA-11 12 13 Time (Min) 08S1-401207F 09S1-701207A-08S1-noloonA 09S1-roloorA Aroclor+1254 -21 . - m . ₽621-roloorA - co ₽621-701207<u>6</u> . -r . . -w analyx-m-proidosataT -ហ 0.38 0.40-

Data File: T1230597.D

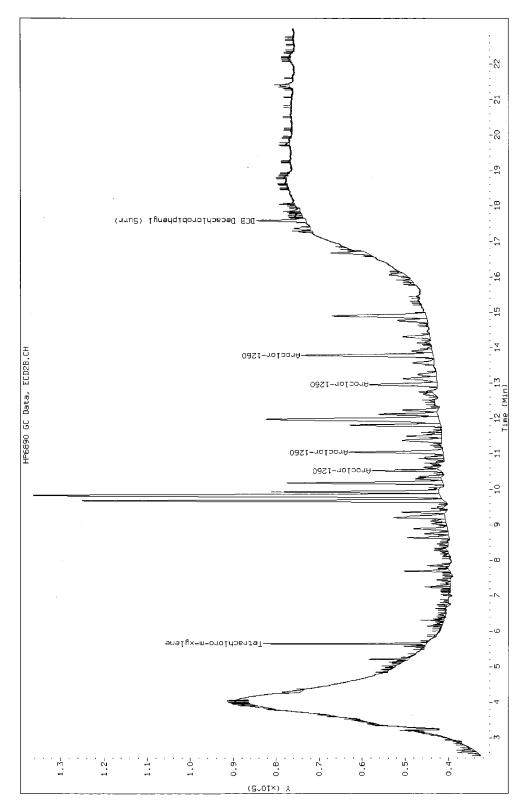
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Date: 28-DEC-2013 05:07

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i



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Data File: T1230598.D

Date: 28-DEC-2013 05:39

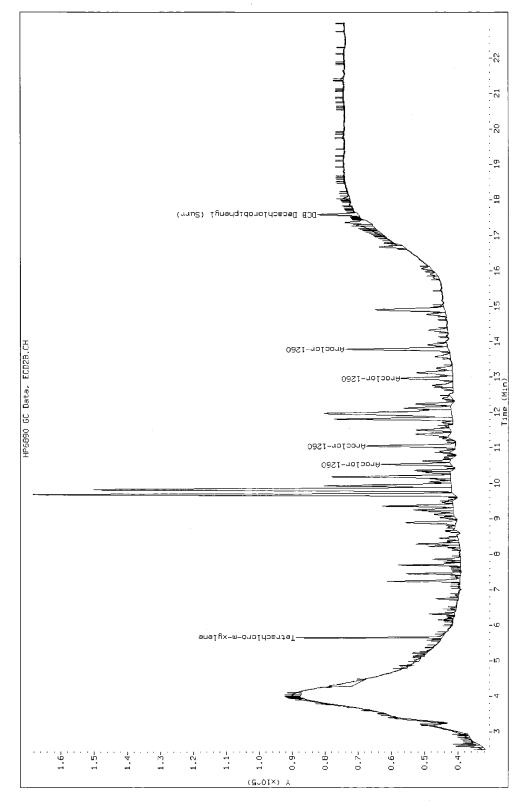
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SD-194-0-1

Operator: 402360

Instrument: gc10.i



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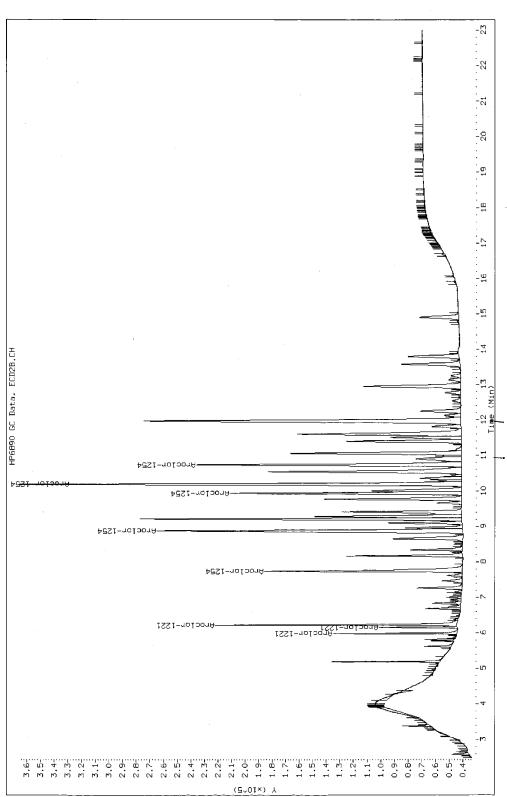
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Date: 27-DEC-2013 00:46

Client ID:

Sample Info: 12273B8082ALL.b

Instrument: gc10.i Operator: 402360 HP6890 GC Data, ECD2B.CH



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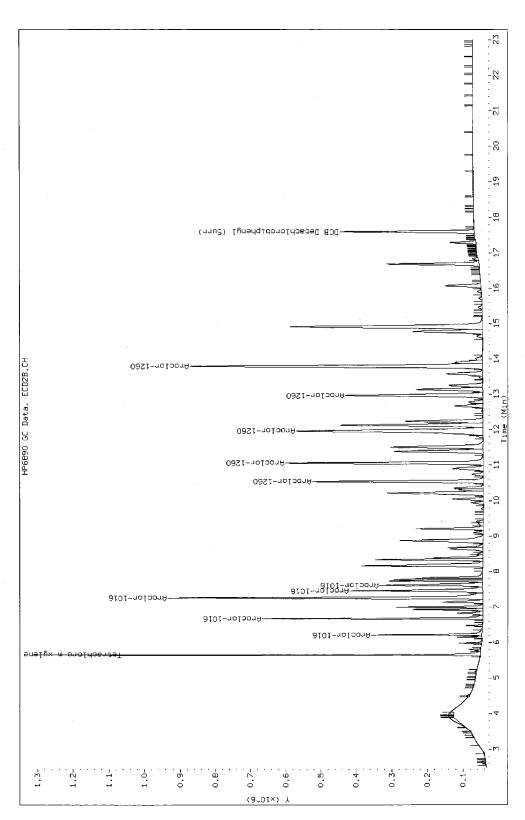
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Instrument: gc10.i



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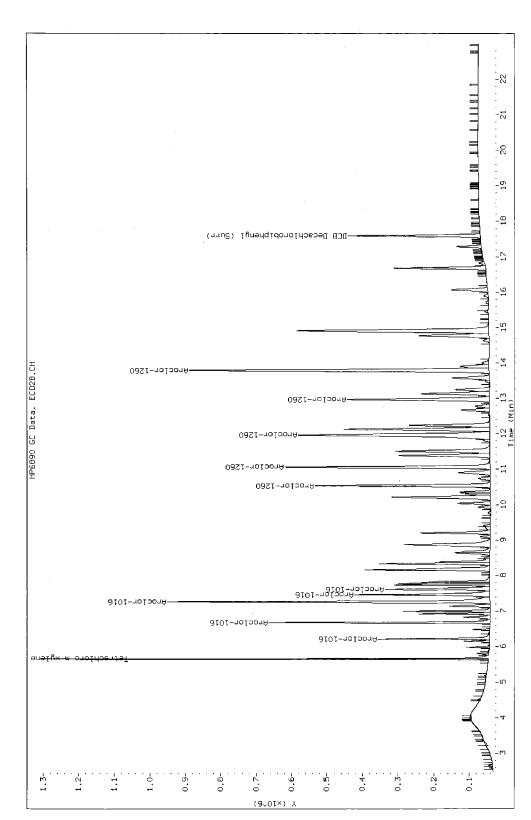
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Sample Info: 12273B8082ALL.b

Instrument: gc10.i



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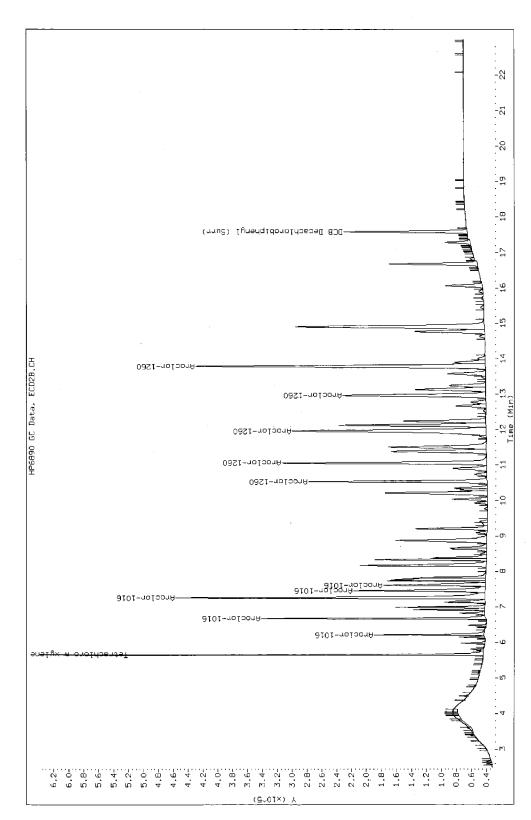
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Sample Info: 12273B8082ALL.b







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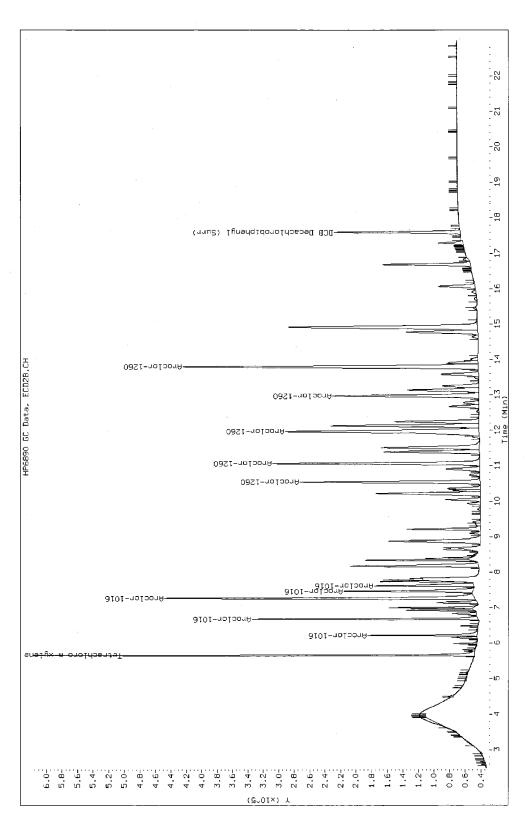
Date: 28-DEC-2013 15:07

Client ID:

Sample Info: 12273B8082ALL.b

BS

Instrument: gc10.i



APPENDIX C—CHEMICAL DATA TABLE									

Chemical Results for Storm Drainage System Sediment Samples-2013

Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland

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LOCATION	SD-163	SD-164	SD-165	SD-166	SD-167	SD-168	SD-169	SD-170
SAMPLE ID	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-167-0-1	SD-168-0-1	SD-169-0-1	SD-170-0-1
SAMPLE DATE	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/20/2013
POLYCYCLIC AROMATIC HYDRO	CARBONS (UG/KG)							
BaPEq-POS	775.31	966.96	1429.4	233.18	382.14	592.83	95.75	362.06
2-METHYLNAPHTHALENE	19 U	30 J	25 J	75 U	39 U	38 U	17 U	19 U
ACENAPHTHENE	64 J	25 J	47 J	80 U	41 U	110 J	30 J	20 U
ACENAPHTHYLENE	69 J	150 J	36 J	96 U	49 U	48 U	22 U	58 J
ANTHRACENE	97 J	200 J	110 J	82 U	44 J	180 J	19 U	110 J
BENZO(A)ANTHRACENE	530 J	520	830	130 J	270 J	470	76 J	320
BENZO(A)PYRENE	500 J	660	920	120 J	280 J	390 J	69 J	250 J
BENZO(B)FLUORANTHENE	700 J	640	1400	130 U	430	250 J	120 J	320 J
BENZO(G,H,I)PERYLENE	480 J	760	890	130 J	340 J	410 J	81 J	170 J
BENZO(K)FLUORANTHENE	260 J	420	590	170 U	170 J	530	43 J	78 J
CHRYSENE	710 J	760	1500	180 J	440	530	120 J	280
DIBENZO(A,H)ANTHRACENE	110 J	130 J	200 J	100 J	48 U	88 J	21 UJ	32 J
FLUORANTHENE	1400 J	960	3500	180 J	800	1000	180 J	630
FLUORENE	52 J	52 J	55 J	110 U	57 U	86 J	25 U	28 U
INDENO(1,2,3-CD)PYRENE	390 J	560	790	86 U	300 J	370 J	66 J	150 J
NAPHTHALENE	19 U	34 J	95 J	72 U	37 U	36 J	16 U	18 U
PHENANTHRENE	990 J	370	2800	130 U	340 J	710	47 J	300
PYRENE	1100 J	1100	2500	210 J	590	830	160 J	480
PCBS (UG/KG)								
AROCLOR-1016	0.4 U	0.39 U	0.41 U	0.39 U	0.4 U	0.39 U	0.35 U	0.39 U
AROCLOR-1221	0.52 U	0.5 U	0.52 U	0.5 U	0.51 U	0.5 U	0.45 U	0.51 U
AROCLOR-1232	0.46 U	0.45 U	0.47 U	0.45 U	0.46 U	0.45 U	0.41 U	0.45 U
AROCLOR-1242	0.44 U	0.43 U	0.44 U	0.42 U	0.44 U	0.43 U	0.39 U	0.43 U
AROCLOR-1248	0.26 U	0.25 U	0.26 U	0.25 U	0.25 U	0.25 U	0.22 U	0.25 U
AROCLOR-1254	0.38 U	0.38 U	0.39 U	0.37 U	0.38 U	100	29	0.38 U
AROCLOR-1260	25	89	99 J	34	19	84	29	2.1 J
TOTAL AROCLOR	25	89	99	34	19	184	58	2.1

Chemical Results for Storm Drainage System Sediment Samples-2013

Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland

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LOCATION	SD-163	SD-164	SD-165	SD-166	SD-171	SD-172	SD-173	SD-174	
SAMPLE ID	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-171-0-1	SD-172-0-3	SD-173-0-3	SD-174-0-1	
SAMPLE DATE	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/19/2013	12/19/2013	12/19/2013	12/19/2013	
POLYCYCLIC AROMATIC HYDROCARBONS (UG/KG)									
BaPEq-POS	775.31	966.96	1429.4	233.18	1020.22	6393	2828.3	2870.2	
2-METHYLNAPHTHALENE	19 U	30 J	25 J	75 U	14 J	50 J	55 J	47 J	
ACENAPHTHENE	64 J	25 J	47 J	80 U	28 J	550	150 J	230	
ACENAPHTHYLENE	69 J	150 J	36 J	96 U	71	160 J	430	140 J	
ANTHRACENE	97 J	200 J	110 J	82 U	160	1100	580	610	
BENZO(A)ANTHRACENE	530 J	520	830	130 J	620	4600	1700	2200	
BENZO(A)PYRENE	500 J	660	920	120 J	680 J	4200 J	1900 J	1900 J	
BENZO(B)FLUORANTHENE	700 J	640	1400	130 U	720 J	5300 J	2100 J	2500 J	
BENZO(G,H,I)PERYLENE	480 J	760	890	130 J	630 J	3300 J	1800 J	1300 J	
BENZO(K)FLUORANTHENE	260 J	420	590	170 U	250 J	1800 J	630 J	780 J	
CHRYSENE	710 J	760	1500	180 J	720	5000	2000	2400	
DIBENZO(A,H)ANTHRACENE	110 J	130 J	200 J	100 J	150 J	870 J	390 J	360 J	
FLUORANTHENE	1400 J	960	3500	180 J	910	9600	3000	5800	
FLUORENE	52 J	52 J	55 J	110 U	35 J	460	160 J	250	
INDENO(1,2,3-CD)PYRENE	390 J	560	790	86 U	530 J	3100 J	1500 J	1300 J	
NAPHTHALENE	19 U	34 J	95 J	72 U	24 J	89 J	94 J	160 J	
PHENANTHRENE	990 J	370	2800	130 U	430	5000	1100	3800	
PYRENE	1100 J	1100	2500	210 J	1100	7400	3400	4000	
PCBS (UG/KG)									
AROCLOR-1016	0.4 U	0.39 U	0.41 U	0.39 U	190 U	120 U	500 U	0.37 U	
AROCLOR-1221	0.52 U	0.5 U	0.52 U	0.5 U	240 U	150 U	640 U	0.48 U	
AROCLOR-1232	0.46 U	0.45 U	0.47 U	0.45 U	210 U	130 U	580 U	0.43 U	
AROCLOR-1242	0.44 U	0.43 U	0.44 U	0.42 U	200 U	130 U	550 U	0.41 U	
AROCLOR-1248	0.26 U	0.25 U	0.26 U	0.25 U	120 U	74 U	320 U	0.24 U	
AROCLOR-1254	0.38 U	0.38 U	0.39 U	0.37 U	180 U	110 U	480 U	150	
AROCLOR-1260	25	89	99 J	34	150000	54000	220000	350	
TOTAL AROCLOR	25	89	99	34	150000	54000	220000	500	

Chemical Results for Storm Drainage System Sediment Samples-2013

Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland

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LOCATION	SD-163	SD-164	SD-165	SD-166	SD-175	SD-176	SD-177	SD-178
SAMPLE ID	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-175-0-1	SD-176-0-3	SD-177-0-2	SD-178-0-1
SAMPLE DATE	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/19/2013	12/19/2013	12/19/2013	12/19/2013
POLYCYCLIC AROMATIC HYDROG	CARBONS (UG/KG)			_				
BaPEq-POS	775.31	966.96	1429.4	233.18	8061.1	2699.3	30526	258.05
2-METHYLNAPHTHALENE	19 U	30 J	25 J	75 U	210 J	30 J	300 J	19 U
ACENAPHTHENE	64 J	25 J	47 J	80 U	1500	130 J	2200 J	20 U
ACENAPHTHYLENE	69 J	150 J	36 J	96 U	310	200 J	510 J	48 J
ANTHRACENE	97 J	200 J	110 J	82 U	2100	390	3300 J	30 J
BENZO(A)ANTHRACENE	530 J	520	830	130 J	5700	1600	18000 J	140 J
BENZO(A)PYRENE	500 J	660	920	120 J	5400 J	1800 J	21000 J	180 J
BENZO(B)FLUORANTHENE	700 J	640	1400	130 U	6000 J	2700 J	27000 J	200 J
BENZO(G,H,I)PERYLENE	480 J	760	890	130 J	3700 J	1500 J	15000 J	140 J
BENZO(K)FLUORANTHENE	260 J	420	590	170 U	2500 J	710 J	10000 J	86 J
CHRYSENE	710 J	760	1500	180 J	6100	2200	26000 J	190 J
DIBENZO(A,H)ANTHRACENE	110 J	130 J	200 J	100 J	1100 J	320 J	3500 J	32 J
FLUORANTHENE	1400 J	960	3500	180 J	13000	4200	68000 J	260
FLUORENE	52 J	52 J	55 J	110 U	1000	160 J	2200 J	28 U
INDENO(1,2,3-CD)PYRENE	390 J	560	790	86 U	3600 J	1400 J	14000 J	110 J
NAPHTHALENE	19 U	34 J	95 J	72 U	480	51 J	120 J	18 U
PHENANTHRENE	990 J	370	2800	130 U	9000	1700	39000 J	81 J
PYRENE	1100 J	1100	2500	210 J	10000	3000	45000 J	210
PCBS (UG/KG)								
AROCLOR-1016	0.4 U	0.39 U	0.41 U	0.39 U	4.4 U	0.55 U	1.1 UJ	0.39 U
AROCLOR-1221	0.52 U	0.5 U	0.52 U	0.5 U	5.7 U	0.7 U	1.4 UJ	0.5 U
AROCLOR-1232	0.46 U	0.45 U	0.47 U	0.45 U	5.1 U	0.63 U	1.3 UJ	0.45 U
AROCLOR-1242	0.44 U	0.43 U	0.44 U	0.42 U	4.8 U	0.6 U	1.2 UJ	0.43 U
AROCLOR-1248	0.26 U	0.25 U	0.26 U	0.25 U	2.8 U	0.35 U	0.69 UJ	0.25 U
AROCLOR-1254	0.38 U	0.38 U	0.39 U	0.37 U	410	36	28 J	5.3
AROCLOR-1260	25	89	99 J	34	1300	88	73 J	12
TOTAL AROCLOR	25	89	99	34	1710	124	101	17.3

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Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland

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LOCATION	SD-163	SD-164	SD-165	SD-166	SD-179	SD-180	SD-181	SD-182
SAMPLE ID	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-179-0-1	SD-180-0-1	SD-181-0-1	SD-182-0-3
SAMPLE DATE	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/19/2013	12/19/2013	12/20/2013	12/20/2013
POLYCYCLIC AROMATIC HYDROC	ARBONS (UG/KG)							
BaPEq-POS	775.31	966.96	1429.4	233.18	289.43	4322.7	25508	14776
2-METHYLNAPHTHALENE	19 U	30 J	25 J	75 U	2.5 J	38 U	440 J	240 J
ACENAPHTHENE	64 J	25 J	47 J	80 U	18	70 J	3900	1600
ACENAPHTHYLENE	69 J	150 J	36 J	96 U	14	76 J	350 J	230 J
ANTHRACENE	97 J	200 J	110 J	82 U	41	390 J	7100	3000
BENZO(A)ANTHRACENE	530 J	520	830	130 J	120	2600	20000	10000
BENZO(A)PYRENE	500 J	660	920	120 J	200 J	2900	17000	9500
BENZO(B)FLUORANTHENE	700 J	640	1400	130 U	290 J	4300	21000	13000
BENZO(G,H,I)PERYLENE	480 J	760	890	130 J	150 J	2600	14000	9200
BENZO(K)FLUORANTHENE	260 J	420	590	170 U	120 J	1900	8800	3500
CHRYSENE	710 J	760	1500	180 J	230	3700	20000	11000
DIBENZO(A,H)ANTHRACENE	110 J	130 J	200 J	100 J	32 J	490	3100	2200
FLUORANTHENE	1400 J	960	3500	180 J	410	6200	35000	18000
FLUORENE	52 J	52 J	55 J	110 U	13	110 J	4100	1200
INDENO(1,2,3-CD)PYRENE	390 J	560	790	86 U	150 J	2200	12000	7300
NAPHTHALENE	19 U	34 J	95 J	72 U	4.7 J	36 U	640	450 J
PHENANTHRENE	990 J	370	2800	130 U	150	2600	28000	12000
PYRENE	1100 J	1100	2500	210 J	290	5900	37000	18000
PCBS (UG/KG)								
AROCLOR-1016	0.4 U	0.39 U	0.41 U	0.39 U	0.44 U	0.79 U	2100 U	2600 U
AROCLOR-1221	0.52 U	0.5 U	0.52 U	0.5 U	0.57 U	1 U	2800 U	3300 U
AROCLOR-1232	0.46 U	0.45 U	0.47 U	0.45 U	0.51 U	0.91 U	2500 U	3000 U
AROCLOR-1242	0.44 U	0.43 U	0.44 U	0.42 U	0.49 U	0.86 U	2400 U	2800 U
AROCLOR-1248	0.26 U	0.25 U	0.26 U	0.25 U	0.28 U	0.5 U	1400 U	1600 U
AROCLOR-1254	0.38 U	0.38 U	0.39 U	0.37 U	7.6	21	2100 U	2500 U
AROCLOR-1260	25	89	99 J	34	22	95	420000	780000
TOTAL AROCLOR	25	89	99	34	29.6	116	420000	780000

Chemical Results for Storm Drainage System Sediment Samples-2013

Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland

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LOCATION	SD-163	SD-164	SD-165	SD-166	SD-183	SD-184	SD-185	SD-186			
SAMPLE ID	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-183-0-1	SD-184-0-2	SD-185-0-1	SD-186-0-1			
SAMPLE DATE	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013			
POLYCYCLIC AROMATIC HYDROC	POLYCYCLIC AROMATIC HYDROCARBONS (UG/KG)										
BaPEq-POS	775.31	966.96	1429.4	233.18	20148	5971.9	1857.3	471.02			
2-METHYLNAPHTHALENE	19 U	30 J	25 J	75 U	280	170	21 J	32 U			
ACENAPHTHENE	64 J	25 J	47 J	80 U	3000	920	53 J	35 U			
ACENAPHTHYLENE	69 J	150 J	36 J	96 U	320	560	94 J	50 J			
ANTHRACENE	97 J	200 J	110 J	82 U	5600	2200	230	96 J			
BENZO(A)ANTHRACENE	530 J	520	830	130 J	15000	4500	1300 J	280 J			
BENZO(A)PYRENE	500 J	660	920	120 J	13000	3800	1200 J	290 J			
BENZO(B)FLUORANTHENE	700 J	640	1400	130 U	15000	4700	1500 J	540			
BENZO(G,H,I)PERYLENE	480 J	760	890	130 J	11000	3400	1100	350 J			
BENZO(K)FLUORANTHENE	260 J	420	590	170 U	6200	1700	570	150 J			
CHRYSENE	710 J	760	1500	180 J	16000	4900	1600	520			
DIBENZO(A,H)ANTHRACENE	110 J	130 J	200 J	100 J	3100	930	270	67 J			
FLUORANTHENE	1400 J	960	3500	180 J	33000	11000	3600	1100			
FLUORENE	52 J	52 J	55 J	110 U	2400	1100	87 J	47 U			
INDENO(1,2,3-CD)PYRENE	390 J	560	790	86 U	9700	3000	1000	300 J			
NAPHTHALENE	19 U	34 J	95 J	72 U	480	340	43 J	31 U			
PHENANTHRENE	990 J	370	2800	130 U	17000	6400	1100 J	400			
PYRENE	1100 J	1100	2500	210 J	19000	6100	1900	590			
PCBS (UG/KG)											
AROCLOR-1016	0.4 U	0.39 U	0.41 U	0.39 U	410 U	0.91 U	0.83 U	0.84 U			
AROCLOR-1221	0.52 U	0.5 U	0.52 U	0.5 U	530 U	1.2 U	1.1 U	1.1 U			
AROCLOR-1232	0.46 U	0.45 U	0.47 U	0.45 U	470 U	1 U	0.96 U	0.96 U			
AROCLOR-1242	0.44 U	0.43 U	0.44 U	0.42 U	450 U	0.99 U	0.91 U	0.91 U			
AROCLOR-1248	0.26 U	0.25 U	0.26 U	0.25 U	260 U	0.58 U	0.53 U	0.53 U			
AROCLOR-1254	0.38 U	0.38 U	0.39 U	0.37 U	390 U	410	190	38			
AROCLOR-1260	25	89	99 J	34	40000	1300	550	60			
TOTAL AROCLOR	25	89	99	34	40000	1710	740	98			

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Lockheed Martin Middle River Complex and Martin State Airport, Middle River, Maryland

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LOCATION	SD-163	SD-164	SD-165	SD-166	SD-187	SD-188	SD-189	SD-190		
SAMPLE ID	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-187-0-1	SD-188-0-1	SD-189-0-1	SD-190-0-1		
SAMPLE DATE	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013		
POLYCYCLIC AROMATIC HYDROC	POLYCYCLIC AROMATIC HYDROCARBONS (UG/KG)									
BaPEq-POS	775.31	966.96	1429.4	233.18	2168.9	954.21	4948.7	64045		
2-METHYLNAPHTHALENE	19 U	30 J	25 J	75 U	48 J	27 U	26 U	530		
ACENAPHTHENE	64 J	25 J	47 J	80 U	48 U	29 U	130 J	5600		
ACENAPHTHYLENE	69 J	150 J	36 J	96 U	190 J	120 J	160 J	870		
ANTHRACENE	97 J	200 J	110 J	82 U	260 J	140 J	590	18000		
BENZO(A)ANTHRACENE	530 J	520	830	130 J	1100	470	3000	49000		
BENZO(A)PYRENE	500 J	660	920	120 J	1400	620	3300	43000		
BENZO(B)FLUORANTHENE	700 J	640	1400	130 U	1800	800	3900	47000		
BENZO(G,H,I)PERYLENE	480 J	760	890	130 J	1600	690	2800	33000		
BENZO(K)FLUORANTHENE	260 J	420	590	170 U	720	250 J	1500	19000		
CHRYSENE	710 J	760	1500	180 J	1700	710	3700	55000		
DIBENZO(A,H)ANTHRACENE	110 J	130 J	200 J	100 J	340 J	150 J	680	8200		
FLUORANTHENE	1400 J	960	3500	180 J	2900	1000	8100	150000		
FLUORENE	52 J	52 J	55 J	110 U	82 J	43 J	180 J	6600		
INDENO(1,2,3-CD)PYRENE	390 J	560	790	86 U	1300	540	2600	30000		
NAPHTHALENE	19 U	34 J	95 J	72 U	43 U	26 U	66 J	120 J		
PHENANTHRENE	990 J	370	2800	130 U	850	260 J	2700	61000		
PYRENE	1100 J	1100	2500	210 J	1900	710	4500	74000		
PCBS (UG/KG)										
AROCLOR-1016	0.4 U	0.39 U	0.41 U	0.39 U	0.78 U	0.7 U	0.68 U	0.88 U		
AROCLOR-1221	0.52 U	0.5 U	0.52 U	0.5 U	0.99 U	0.9 U	0.87 U	1.1 U		
AROCLOR-1232	0.46 U	0.45 U	0.47 U	0.45 U	0.89 U	0.81 U	0.78 U	1 U		
AROCLOR-1242	0.44 U	0.43 U	0.44 U	0.42 U	0.85 U	0.77 U	0.75 U	0.97 U		
AROCLOR-1248	0.26 U	0.25 U	0.26 U	0.25 U	0.49 U	0.45 U	0.43 U	0.56 U		
AROCLOR-1254	0.38 U	0.38 U	0.39 U	0.37 U	29	15	11	53		
AROCLOR-1260	25	89	99 J	34	92	49	82	46		
TOTAL AROCLOR	25	89	99	34	121	64	93	99		

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LOCATION	SD-163	SD-164	SD-165	SD-166	SD-191	SD-192	SD-193	SD-194		
SAMPLE ID	SD-163-0-1	SD-164-0-1	SD-165-0-1	SD-166-0-1	SD-191-0-1	SD-192-0-1	SD-193-0-1	SD-194-0-1		
SAMPLE DATE	12/18/2013	12/18/2013	12/18/2013	12/18/2013	12/20/2013	12/20/2013	12/20/2013	12/20/2013		
POLYCYCLIC AROMATIC HYDROCARBONS (UG/KG)										
BaPEq-POS	775.31	966.96	1429.4	233.18	19123	219.73	251.07	188.49		
2-METHYLNAPHTHALENE	19 U	30 J	25 J	75 U	230	7.8 U	8.4 U	7.1 U		
ACENAPHTHENE	64 J	25 J	47 J	80 U	2000	20 J	13 J	9 J		
ACENAPHTHYLENE	69 J	150 J	36 J	96 U	440	89	31 J	48 J		
ANTHRACENE	97 J	200 J	110 J	82 U	4200	89	54 J	43 J		
BENZO(A)ANTHRACENE	530 J	520	830	130 J	13000	130	160	100		
BENZO(A)PYRENE	500 J	660	920	120 J	13000	150	170	130		
BENZO(B)FLUORANTHENE	700 J	640	1400	130 U	16000	170	200	140		
BENZO(G,H,I)PERYLENE	480 J	760	890	130 J	9200	120	130	94		
BENZO(K)FLUORANTHENE	260 J	420	590	170 U	5600	64 J	87 J	44 J		
CHRYSENE	710 J	760	1500	180 J	17000	190	200	150		
DIBENZO(A,H)ANTHRACENE	110 J	130 J	200 J	100 J	2300	29 J	33 J	26 J		
FLUORANTHENE	1400 J	960	3500	180 J	67000	360	440	240		
FLUORENE	52 J	52 J	55 J	110 U	2400	34 J	23 J	15 J		
INDENO(1,2,3-CD)PYRENE	390 J	560	790	86 U	8500	99	110	79 J		
NAPHTHALENE	19 U	34 J	95 J	72 U	87 J	7.5 U	8.1 U	6.8 U		
PHENANTHRENE	990 J	370	2800	130 U	42000	170	160	81		
PYRENE	1100 J	1100	2500	210 J	49000	240	230	160		
PCBS (UG/KG)										
AROCLOR-1016	0.4 U	0.39 U	0.41 U	0.39 U	0.89 U	0.81 U	0.88 U	0.73 U		
AROCLOR-1221	0.52 U	0.5 U	0.52 U	0.5 U	1.1 U	1 U	1.1 U	0.94 U		
AROCLOR-1232	0.46 U	0.45 U	0.47 U	0.45 U	1 U	0.94 U	1 U	0.84 U		
AROCLOR-1242	0.44 U	0.43 U	0.44 U	0.42 U	0.97 U	0.89 U	0.96 U	0.8 U		
AROCLOR-1248	0.26 U	0.25 U	0.26 U	0.25 U	0.56 U	0.52 U	0.56 U	0.47 U		
AROCLOR-1254	0.38 U	0.38 U	0.39 U	0.37 U	350	13	0.84 U	0.7 U		
AROCLOR-1260	25	89	99 J	34	530	20	17	16		
TOTAL AROCLOR	25	89	99	34	880	33	17	16		

BaPEq - benzo(a)pyrene equivalent

J - estimated concentration

PCBs - polychlorinated biphenyls

POS - only detected polycyclic aromatic hydrocarbons are used for this calculation

U - not detected at the concentration shown left of the letter.

UG/KG - micrograms per kilogram (i.e., parts per billion)

-- not analyzed